

## Optimizing smallpt

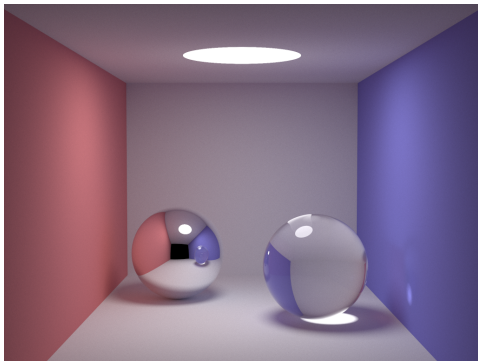
Davean Scies, Siddharth Bhat

## Haskell Exchange

November 4th, 2020



# What is smallpt anyway?

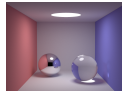


- ▶ 99 LoC C++ raytracer.
- ▶ Perfect for an optimization case study.
- ▶ Ported to many languages, including Haskell! (Thanks to Vo Minh Thu(noteed)).
- ▶ Start from noteed's original source; SHA the output image for baseline, keep optimizing.
- ▶ Plan: Quick walk through Haskell code, end up at C++ (clang++) performance.

## Optimizing smallpt

2020-11-04

What is smallpt anyway?



- 99 LoC C++ raytracer.
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- Plan: Quick walk through Haskell code, end up at C++ (clang++) performance.

Note for first slide: what is smallpt anyway?

# What is smallpt anyway?

```
struct Vec {
    double x, y, z; // position, also color (r,g,b)
    ... methods...
};

struct Ray { Vec o, d; Ray(Vec o_, Vec d_) : o(o_), d(d_) {} };

enum Refl_t { DIFF, SPEC, REFR }; // material types, used in radiance()

struct Sphere {
    double rad; // radius
    Vec p, e, c; // position, emission, color
    Refl_t refl; // reflection type (DIFFuse, SPECular, REFRactive)
    ... methods ...
    double intersect(const Ray &r) const // returns distance, 0 if nohit
};

Sphere spheres[] = { //Scene: radius, position, emission, color, material
    Sphere(1e5, Vec( 1e5+1,40.8,81.6), Vec(),Vec(.75,.25,.25),DIFF), //Left
    ... initialization ...
};
```

## Optimizing smallpt

2020-11-04

What is smallpt anyway?

```
What is smallpt anyway?

struct Vec {
    double x, y, z; // position, also color (r,g,b)
    ... methods...
};

struct Ray { Vec o, d; Ray(Vec o_, Vec d_) : o(o_), d(d_) {} };

enum Refl_t { DIFF, SPEC, REFR }; // material types, used in radiance()

struct Sphere {
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    ... methods ...
    double intersect(const Ray &r) const // returns distance, 0 if nohit
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Sphere spheres[] = { //Scene: radius, position, emission, color, material
    Sphere(1e5, Vec( 1e5+1,40.8,81.6), Vec(),Vec(.75,.25,.25),DIFF), //Left
    ... initialization ...
};
```

# What is smallpt anyway?

```
Vec radiance(const Ray &r, int depth, unsigned short *Xi){
```

```
}
```

## Optimizing smallpt

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What is smallpt anyway?

Say that the core function is radiance



# What is smallpt anyway?

```
Vec radiance(const Ray &r, int depth, unsigned short *Xi){
```

```
    radiance
```

```
    radiance
```

```
    radiance
```

```
    radiance  
    radiance
```

```
    radiance  
    radiance
```

```
}
```

## Optimizing smallpt

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What is smallpt anyway?



# What is smallpt anyway?

```
Vec radiance(const Ray &r, int depth, unsigned short *Xi){
```

```
    if ( ) if ( ) else  
    if ( ){
```

```
        radiance  
    } else if ( )  
        radiance
```

```
    if ( )  
        radiance
```

```
        radiance  
        radiance  
    }  
}
```

## Optimizing smallpt

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What is smallpt anyway?

...with control flow

```
What is smallpt anyway?  
  
Vec radiance(const Ray &r, int depth, unsigned short *Xi){  
  
    if ( ) if ( ) else  
    if ( ){  
  
        radiance  
    } else if ( )  
        radiance  
  
    if ( )  
        radiance  
  
        radiance  
        radiance  
    }  
}
```

# What is smallpt anyway?

```
Vec radiance(const Ray &r, int depth, unsigned short *Xi){
```

```
Vec x=r.o+r.d*t, n=(x-obj.p).norm(), nl=n.dot(r.d)<0?n:n*-1, f=obj.c;
```

```
if ( ) if ( ) else  
if ( ){
```

```
radiance  
} else if ( )  
radiance
```

```
if ((cos2t=1-nnt*nnt*(1-ddn*ddn))<0)  
radiance
```

```
radiance radiance  
radiance radiance  
}
```

## Optimizing smallpt

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What is smallpt anyway?

and lots of arithmetic

```
What is smallpt anyway?  
  
Vec radiance(const Ray &r, int depth, unsigned short *Xi){  
  
Vec w=obj.w, w1=obj.g*w, w2=obj.b*w, w3=obj.r*w, w4=obj.y;  
if ( ) if ( ) else  
if ( ){  
  
radiance  
radiance  
}  
else if ( )  
radiance  
radiance  
}  
if ((cos2t=1-nnt*nnt*(1-ddn*ddn))<0)  
radiance  
  
radiance radiance  
radiance radiance  
}
```



# What is smallpt anyway?

```
Vec radiance(const Ray &r, int depth, unsigned short *Xi){
    double t; // distance to intersection
    int id=0; // id of intersected object
    if (!intersect(r, t, id)) return Vec(); // if miss, return black
    const Sphere &obj = spheres[id]; // the hit object
    Vec x=r.o+r.d*t, n=(x-obj.p).norm(), nl=n.dot(r.d)<0?n:n*-1, f=obj.c;
    double p = f.x>f.y && f.x>f.z ? f.x : f.y>f.z ? f.y : f.z; // max refl
    if (++depth>5) if (erand48(Xi)<p) f=f*(1/p); else return obj.e; //R.R.
    if (obj.refl == DIFF){ // Ideal DIFFUSE reflection
        double r1=2*M_PI*erand48(Xi), r2=erand48(Xi), r2s=sqrt(r2);
        Vec w=nl, u=((fabs(w.x)>.1?Vec(0,1):Vec(1))%w).norm(), v=w%u;
        Vec d = (u*cos(r1)*r2s + v*sin(r1)*r2s + w*sqrt(1-r2)).norm();
        return obj.e + f.mult(radiance(Ray(x,d),depth,Xi));
    } else if (obj.refl == SPEC) // Ideal SPECULAR reflection
        return obj.e + f.mult(radiance(Ray(x,r.d-n*2*n.dot(r.d)),depth,Xi));
    Ray reflRay(x, r.d-n*2*n.dot(r.d)); // Ideal dielectric REFRACTION
    bool into = n.dot(nl)>0; // Ray from outside going in?
    double nc=1, nt=1.5, nnt=into?nc/nt:nt/nc, ddn=r.d.dot(nl), cos2t;
    if ((cos2t=1-nnt*nnt*(1-ddn*ddn))<0) // Total internal reflection
        return obj.e + f.mult(radiance(reflRay,depth,Xi));
    Vec tdir = (r.d*nnt - n*((into?-1):1)*(ddn*nnt+sqrt(cos2t))).norm();
    double a=nt-nc, b=nt+nc, R0=a*a/(b*b), c = 1-(into?-ddn:tdir.dot(n));
    double Re=R0+(1-R0)*c*c*c*c*c, Tr=1-Re, P=.25+.5*Re, RP=Re/P, TP=Tr/(1-P);
    return obj.e + f.mult(depth>2 ? (erand48(Xi)<P ? // Russian roulette
        radiance(reflRay,depth,Xi)*RP:radiance(Ray(x,tdir),depth,Xi)*TP) :
        radiance(reflRay,depth,Xi)*Re+radiance(Ray(x,tdir),depth,Xi)*Tr);
}
```

## Optimizing smallpt

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What is smallpt anyway?

display the whole thing

What is smallpt anyway?

```
Vec radiance(const Ray &r, int depth, unsigned short *Xi){
    double t; // distance to intersection
    int id=0; // id of intersected object
    if (!intersect(r, t, id)) return Vec(); // if miss, return black
    const Sphere &obj = spheres[id]; // the hit object
    Vec x=r.o+r.d*t, n=(x-obj.p).norm(), nl=n.dot(r.d)<0?n:n*-1, f=obj.c;
    double p = f.x>f.y && f.x>f.z ? f.x : f.y>f.z ? f.y : f.z; // max refl
    if (++depth>5) if (erand48(Xi)<p) f=f*(1/p); else return obj.e; //R.R.
    if (obj.refl == DIFF){ // Ideal DIFFUSE reflection
        double r1=2*M_PI*erand48(Xi), r2=erand48(Xi), r2s=sqrt(r2);
        Vec w=nl, u=((fabs(w.x)>.1?Vec(0,1):Vec(1))%w).norm(), v=w%u;
        Vec d = (u*cos(r1)*r2s + v*sin(r1)*r2s + w*sqrt(1-r2)).norm();
        return obj.e + f.mult(radiance(Ray(x,d),depth,Xi));
    } else if (obj.refl == SPEC) // Ideal SPECULAR reflection
        return obj.e + f.mult(radiance(Ray(x,r.d-n*2*n.dot(r.d)),depth,Xi));
    Ray reflRay(x, r.d-n*2*n.dot(r.d)); // Ideal dielectric REFRACTION
    bool into = n.dot(nl)>0; // Ray from outside going in?
    double nc=1, nt=1.5, nnt=into?nc/nt:nt/nc, ddn=r.d.dot(nl), cos2t;
    if ((cos2t=1-nnt*nnt*(1-ddn*ddn))<0) // Total internal reflection
        return obj.e + f.mult(radiance(reflRay,depth,Xi));
    Vec tdir = (r.d*nnt - n*((into?-1):1)*(ddn*nnt+sqrt(cos2t))).norm();
    double a=nt-nc, b=nt+nc, R0=a*a/(b*b), c = 1-(into?-ddn:tdir.dot(n));
    double Re=R0+(1-R0)*c*c*c*c*c, Tr=1-Re, P=.25+.5*Re, RP=Re/P, TP=Tr/(1-P);
    return obj.e + f.mult(depth>2 ? (erand48(Xi)<P ? // Russian roulette
        radiance(reflRay,depth,Xi)*RP:radiance(Ray(x,tdir),depth,Xi)*TP) :
        radiance(reflRay,depth,Xi)*Re+radiance(Ray(x,tdir),depth,Xi)*Tr);
}
```

# Initial Haskell Code: radiance (1×)

```
radiance :: Ray -> CInt -> Ptr CUSHort -> IO Vec
radiance ray@(Ray o d) depth xi = case intersects ray of
  (Nothing,_) -> return zerov
  (Just t,Sphere _r p e c refl) -> do
```

```
  continue f = case refl of
    DIFF -> do
```

```
      radiance
```

```
    SPEC -> do
```

```
      rad <- radiance
```

```
    REFR -> do
```

```
      if
      then do
        rad <- radiance reflRay depth' xi
```

## Optimizing smallpt

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### Initial Haskell Code: radiance (1×)

show the same thing, this time in haskell

Initial Haskell Code: radiance (1×)

```
radiance :: Ray -> CInt -> Ptr CUSHort -> IO Vec
radiance ray@(Ray o d) depth xi = case intersects ray of
  (Nothing,_) -> return zerov
  (Just t,Sphere _r p e c refl) -> do
```

```
    continue f = case refl of
```

```
      DIFF -> do
```

```
          radiance
```

```
        SPEC -> do
```

```
          rad <- radiance
```

```
        REFR -> do
```

```
          if
```

```
          then do
```

```
            rad <- radiance reflRay depth' xi
```

# Initial Haskell Code: radiance (1×)

```
radiance :: Ray -> CInt -> Ptr CUSHort -> IO Vec
radiance ray@(Ray o d) depth xi = case intersects ray of
  (Nothing,_) -> return zeroV
  (Just t,Sphere _r p e c refl) -> do
    let x = o `addv` (d `mulvs` t)
        n = norm $ x `subv` p
        nl = if n `dot` d < 0 then n else n `mulvs` (-1)
        pr = maxv c
        depth' = depth + 1
    continue f = case refl of
      DIFF -> do
        r1 <- ((2*pi)* `fmap` erand48 xi)
        r2 <- erand48 xi
        let r2s = sqrt r2
            w@(Vec wx _ _) = nl
            u = norm $ (if abs wx > 0.1 then (Vec 0 1 0) else (Vec 1 0 0)) `cross` w
            v = w `cross` u
            d' = norm $ (u`mulvs`(cos r1*r2s)) `addv` (v`mulvs`(sin r1*r2s)) `addv` (w`mulvs`sqrt (1-r2))
        rad <- radiance (Ray x d') depth' xi
        return $ e `addv` (f `mulv` rad)
      SPEC -> do
        let d' = d `subv` (n `mulvs` (2 * (n`dot`d)))
        rad <- radiance (Ray x d') depth' xi
        return $ e `addv` (f `mulv` rad)
      REFR -> do
        let reflRay = Ray x (d `subv` (n `mulvs` (2* n`dot`d)))
            into = n`dot`nl > 0
            nc = 1
            nt = 1.5
            nnt = if into then nc/nt else nt/nc
            ddn= d`dot`nl
            cos2t = 1-nnt*nnt*(1-ddn*ddn)
        if cos2t<0
          then do
            rad <- radiance reflRay depth' xi
```

## Optimizing smallpt

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### Initial Haskell Code: radiance (1×)

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  (Nothing,_) -> return zeroV
  (Just t,Sphere _r p e c refl) -> do
    let x = o `addv` (d `mulvs` t)
        n = norm $ x `subv` p
        nl = if n `dot` d < 0 then n else n `mulvs` (-1)
        pr = maxv c
        depth' = depth + 1
    continue f = case refl of
      DIFF -> do
        r1 <- ((2*pi)* `fmap` erand48 xi)
        r2 <- erand48 xi
        let r2s = sqrt r2
            w@(Vec wx _ _) = nl
            u = norm $ (if abs wx > 0.1 then (Vec 0 1 0) else (Vec 1 0 0)) `cross` w
            v = w `cross` u
            d' = norm $ (u`mulvs`(cos r1*r2s)) `addv` (v`mulvs`(sin r1*r2s)) `addv` (w`mulvs`sqrt (1-r2))
        rad <- radiance (Ray x d') depth' xi
        return $ e `addv` (f `mulv` rad)
      SPEC -> do
        let d' = d `subv` (n `mulvs` (2 * (n`dot`d)))
        rad <- radiance (Ray x d') depth' xi
        return $ e `addv` (f `mulv` rad)
      REFR -> do
        let reflRay = Ray x (d `subv` (n `mulvs` (2* n`dot`d)))
            into = n`dot`nl > 0
            nc = 1
            nt = 1.5
            nnt = if into then nc/nt else nt/nc
            ddn= d`dot`nl
            cos2t = 1-nnt*nnt*(1-ddn*ddn)
        if cos2t<0
          then do
            rad <- radiance reflRay depth' xi
```

full code

# Initial Haskell Code: Entry point (1×)

```
smallpt :: Int -> Int -> Int -> IO ()
```

```
smallpt w h nsamps = do
```

```
    * * *
    c <- VM.replicate (w * h) 0
    allocArray 3 \xi -> -- Create mutable memory
      flip mapM_ [0..h-1] $ \y -> do -- Loop
        writeXi xi y
        for_ [0..w-1] \x -> do -- Loop
          let i = (h-y-1) * w + x
          for_ [0..1] \sy -> do -- Loop
            for_ [0..1] \sx -> do -- Loop
              r <- newIORef 0 -- Create mutable memory
              for_ [0..samps-1] \s -> do -- Loops, Loops
                r1 <- (2*) <$> erand48 xi
                * * *
                rad <- radiance (Ray (org+d.*140) (norm d)) 0 xi -- Crunch
                * * *
                modifyIORef r (+ rad .* recip (fromIntegral samps)) -- Write
              ci <- VM.unsafeRead c i
              Vec rr rg rb <- readIORef r
              VM.unsafeWrite c i $
                ci + Vec (clamp rr) (clamp rg) (clamp rb) .* 0.25 -- Write
```

## Optimizing smallpt

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### Initial Haskell Code: Entry point (1×)

```
Initial Haskell Code: Entry point (1×)

smallpt :: Int -> Int -> Int -> IO ()
smallpt w h nsamps = do
    * * *
    c <- VM.replicate (w * h) 0
    allocArray 3 \xi -> -- Create mutable memory
      flip mapM_ [0..h-1] $ \y -> do -- Loop
        writeXi xi y
        for_ [0..w-1] \x -> do -- Loop
          let i = (h-y-1) * w + x
          for_ [0..1] \sy -> do -- Loop
            for_ [0..1] \sx -> do -- Loop
              r <- newIORef 0 -- Create mutable memory
              for_ [0..samps-1] \s -> do -- Loops, Loops
                r1 <- (2*) <$> erand48 xi
                * * *
                rad <- radiance (Ray (org+d.*140) (norm d)) 0 xi -- Crunch
                * * *
                modifyIORef r (+ rad .* recip (fromIntegral samps)) -- Write
              ci <- VM.unsafeRead c i
              Vec rr rg rb <- readIORef r
              VM.unsafeWrite c i $
                ci + Vec (clamp rr) (clamp rg) (clamp rb) .* 0.25 -- Write
```

## Initial Haskell Code: File I/O (1×)

```
withFile "image.ppm" WriteMode $ \hdl -> do
  hPrintf hdl "P3\n%d %d\n%d\n" w h (255::Int)
  flip mapM_ [0..w*h-1] \i -> do
    Vec r g b <- VM.unsafeRead c i
    hPrintf hdl "%d %d %d " (toInt r) (toInt g) (toInt b)
```

## Optimizing smallpt

2020-11-04

### Initial Haskell Code: File I/O (1×)

```
Initial Haskell Code: File I/O (1×)

withFile "image.ppm" WriteMode $ \hdl -> do
  hPrintf hdl "P3\n%d %d\n%d\n" w h (255::Int)
  flip mapM_ [0..w*h-1] \i -> do
    Vec r g b <- VM.unsafeRead c i
    hPrintf hdl "%d %d %d " (toInt r) (toInt g) (toInt b)
```

## Initial Haskell Code: RNG (1×)

```
foreign import ccall unsafe "erand48"  
  erand48 :: Ptr CUShort -> IO Double
```

## Optimizing smallpt

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Initial Haskell Code: RNG (1×)

Initial Haskell Code: RNG (1×)

```
foreign import ccall unsafe "erand48"  
  erand48 :: Ptr CUShort -> IO Double
```

## Restrict export list to 'main' (1.13×)

```
-module Main where  
+module Main (main) where
```

## Optimizing smallpt

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Restrict export list to 'main' (1.13×)

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```
-module Main where  
+module Main (main) where
```

## Restrict export list to 'main' (1.13×)

```
-module Main where  
+module Main (main) where
```

- ▶ Exported functions could be used by something unknown.
- ▶ original versions must be available.

## Restrict export list to 'main' (1.13×)

```
-module Main where  
+module Main (main) where  
▶ Exported functions could be used by something unknown.  
▶ original versions must be available.
```



### Restrict export list to 'main' (1.13x)

```
-module Main where
+module Main (main) where
```

- ▶ Exported functions could be used by something unknown.
- ▶ original versions must be available.
- ▶ Make many optimizations unsound.

## Optimizing smallpt

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- Restrict export list to 'main' (1.13×)

Restrict export list to 'main' (1.13x)

- ▶ Exported functions could be used by something unknown
- ▶ original versions must be available.
- ▶ Make many optimizations unound.

The very first thing to do is to let the compiler actually optimize. Exported functions could be used by something unknown to the compiler and thus their original versions must be available. This makes many optimizations look bad or be unreasonable. Explicit export lists tell the compiler what you care about. They aren't just about encapsulation.

Mark entries of Ray and Sphere as UNPACK and Strict ( $1.07\times$ )

```
data Vec = Vec {-# UNPACK #-} !Double
            {-# UNPACK #-} !Double
            {-# UNPACK #-} !Double
```

```
-data Ray = Ray Vec Vec -- origin, direction
+data Ray = Ray {-# UNPACK #-} !Vec {-# UNPACK #-} !Vec -- origin, direction
```

```
data Refl = DIFF | SPEC | REFR -- material types, used in radiance
```

```
-- radius, position, emission, color, reflection
-data Sphere = Sphere Double Vec Vec Vec !Refl
+data Sphere = Sphere {-# UNPACK #-} !Double
+                        {-# UNPACK #-} !Vec
+                        {-# UNPACK #-} !Vec
+                        {-# UNPACK #-} !Vec !Refl
```

```
struct Vec { double x, y, z; }
struct Ray { std::function<Vec()> v; std::function<Vec()> w; };
struct RayUnpack { double xv, yv, int zv;
                  double xw, yw, zw; };

```

## Optimizing smallpt

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Mark entries of Ray and Sphere as UNPACK and Strict (1.07x)

```

Mark entries of Ray and Sphere as UNPACK and Strict (1.07)

static Vec = Vec { < UNPACK 2 > | Vecable
                  < < UNPACK 2 > | Vecable
                  < < UNPACK 2 > | Vecable
                  < < UNPACK 2 > | Vecable }

<data Ray = Ray Vec < UNPACK 2 > | origin, direction
<data Ray = Ray < < UNPACK 2 > < | direction < < UNPACK 2 > > | Vec = < origin, direction
<data Ray1 = RAYP < UNPACK 2 > | RAYPS = material types, used in collidance

< | radius, position, material, color, refraction
<data Sphere = Sphere < UNPACK 2 > | Vec = Vec {Vec1
<data Sphere = Sphere < UNPACK 2 > | Vec = Vec {Vec1
+ < < UNPACK 2 > < | Vec =
+ < < UNPACK 2 > < | Vec =
+ < < UNPACK 2 > < | Vec = Vec1
+ < < UNPACK 2 > < | Vec = Vec1

struct Vec { double x, y, z }
struct Ray { Vec <id function Vec() <id > + function Vec() <id > }
struct RaySphere { double w, pos, col, w;
                  double w, pos, col, w;

```

Strict means that they're evaluated ahead of time, not lazily of course. This can save a bit in their evaluation cost but means we always pay it. It is a trade off but here we know it will be a good one because we have almost perfect demand. Unpacking is different. It removes indirection of doing a memory lookup for components but means we have to copy everything into the data structure that it is unpacked into. We don't unpack ray because we'll be doing a lot of calculations on it and we want the Vec math to fuse and not get copied, when we use the Vecs they should be hot in cache anyway. We do inline Sphere because we statically create them and while they'll also be hot in cache there is no benefit to taking any extra cost at all.

## Optimizing smallpt

### Use a pattern synonym to unpack Refl in Sphere (1.07x)

[illegible]

Until the recent addition of unboxed sums in GHC there was no way to unpack a type with multiple constructors. While unboxed sums exist now, their syntax is quite unpleasant. We're using an older trick to fake the unboxing here instead. In this case it isn't much of a win, but it illustrates the technique.

## Change from maximum on a list to max (1.08×)

```
-maxv (Vec a b c) = maximum [a,b,c]
+maxv (Vec a b c) = max a (max b c)

    let x = o `addv` (d `mulvs` t)
        n = norm $ x `subv` p
        nl = if n `dot` d < 0 then n else n `mulvs` (-1)
-    pr = maxv c
    depth' = depth + 1
    continue f = case refl of
        DIFF -> do
...
    if depth'>5
    then do
        er <- erand48 xi
+    let !pr = maxv c
```

## Optimizing smallpt

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Change from maximum on a list to max (1.08×)

```
Change from maximum on a list to max (1.08×)

-maxv (Vec a b c) = maximum [a,b,c]
+maxv (Vec a b c) = max a (max b c)

    let x = o `addv` (d `mulvs` t)
        n = norm $ x `subv` p
        nl = if n `dot` d < 0 then n else n `mulvs` (-1)
-    pr = maxv c
    depth' = depth + 1
    continue f = case refl of
        DIFF -> do
...
    if depth'>5
    then do
        er <- erand48 xi
+    let !pr = maxv c
```

finicky optimization. GHC does not evaluate at compile time, making optimizations like these necessary

# Convert erand48 to pure Haskell (1.09×)

```
-radiance :: Ray -> CInt -> Ptr CUShort -> IO Vec
+radiance :: Ray -> Int -> IORef Word64 -> IO Vec
radiance ray@(Ray o d) depth xi = case intersects ray of
  (Nothing,_) -> return zeroV
  (Just t,Sphere _r p e c refl) -> do
@@ -153,9 +153,8 @@ smallpt w h nsamps = do
    cx = Vec (fromIntegral w * 0.5135 / fromIntegral h) 0 0
    cy = norm (cx `cross` dir) `mulvs` 0.5135
    c <- VM.replicate (w * h) zeroV
-   allocaArray 3 $ \xi ->
-     flip mapM_ [0..h-1] $ \y -> do
+   xi <- newIORef 0
+   flip mapM_ [0..h-1] $ \y -> do
      writeXi xi y
```

## Optimizing smallpt

2020-11-04

### Convert erand48 to pure Haskell (1.09×)

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```
radiance :: Ray -> CInt -> Ptr CUShort -> IO Vec
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    cy = norm (cx `cross` dir) `mulvs` 0.5135
    c <- VM.replicate (w * h) zeroV
-   allocaArray 3 $ \xi ->
-     flip mapM_ [0..h-1] $ \y -> do
+   xi <- newIORef 0
+   flip mapM_ [0..h-1] $ \y -> do
      writeXi xi y
```

This function isn't used much so it doesn't really impact our performance but often building unnecessary data structures is detrimental.

# Remove mutability: Erand48 Monad

```
-erand48 :: IORef Word64 -> IO Double
-erand48 !t = do
-  r <- readIORef t
+data ET a = ET !Word64 !a deriving Functor
+newtype Erand48 a = Erand48 { runErand48' :: Word64 -> ET a } deriving Functor
+instance Applicative Erand48 where
+instance Monad Erand48 where
+runWithErand48 :: Int -> Erand48 a -> a
+erand48 :: Erand48 Double
+...
-radiance :: Ray -> Int -> IORef Word64 -> IO Vec
-radiance ray@(Ray o d) depth xi = case intersects ray of
+radiance :: Ray -> Int -> Erand48 Vec
+radiance ray@(Ray o d) depth = case intersects ray of
+...
-      r1 <- (2*pi*) <$> erand48 xi
-      r2 <- erand48 xi
+      r1 <- (2*pi*) <$> erand48
+      r2 <- erand48
+...
-      then (* rp) <$> radiance reflRay depth' xi
-      else (* tp) <$> radiance (Ray x tdir) depth' xi
+      then (* rp) <$> radiance reflRay depth'
+      else (* tp) <$> radiance (Ray x tdir) depth'
```

# Optimizing smallpt

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## Remove mutability: Erand48 Monad

The entire premise of this talk is that Haskell can be as fast as C. Further any impedance mismatch, such as FFI almost universally has to have, carries some bookkeeping overhead. If our Haskell code was as fast as the C code moving the code into Haskell would be a win, if it was slightly slower it could still be a win. Often considering your Haskell code's performance is a better option and easier than reimplementing something in C. As is the way with optimizations, this is not universally true.

```
Remove mutability: Erand48 Monad
-- Erand48 :: IORef Word64 -> IO Double
-- erand48 !t = do
--   r <- readIORef t
+data ET a = ET !Word64 !a deriving Functor
+newtype Erand48 a = Erand48 { runErand48' :: Word64 -> ET a } deriving Functor
+instance Applicative Erand48 where
+instance Monad Erand48 where
+runWithErand48 :: Int -> Erand48 a -> a
+erand48 :: Erand48 Double
+...
-radiance :: Ray -> Int -> IORef Word64 -> IO Vec
-radiance ray@(Ray o d) depth xi = case intersects ray of
+radiance :: Ray -> Int -> Erand48 Vec
+radiance ray@(Ray o d) depth = case intersects ray of
+...
-      r1 <- (2*pi*) <$> erand48 xi
-      r2 <- erand48 xi
+      r1 <- (2*pi*) <$> erand48
+      r2 <- erand48
+...
-      then (* rp) <$> radiance reflRay depth' xi
-      else (* tp) <$> radiance (Ray x tdir) depth' xi
+      then (* rp) <$> radiance reflRay depth'
+      else (* tp) <$> radiance (Ray x tdir) depth'
```

## Removing mutation: eliminate IORef

```
- c <- VM.replicate (w * h) 0
- xi <- newIORef 0
- flip mapM_ [0..h-1] $ \y -> do
-   writeXi xi y
-   for_ [0..w-1] \x -> do
-     let i = (h-y-1) * w + x
-     for_ [0..1] \sy -> do
-       for_ [0..1] \sx -> do
-         r <- newIORef 0
-         for_ [0..samps-1] \_s -> do
-           r1 <- (2*) <$> erand48 xi
+ img = (`concatMap` [(h-1),(h-2)..0]) $ \y -> runWithErand48 y do
+   for [0..w-1] \x -> do
+     (\pf -> foldlM pf 0 [(sy, sx) | sy <- [0,1], sx <- [0,1]]) \ci (sy, sx) -> do
+       Vec rr rg rb <- (\f -> foldlM f 0 [0..samps-1]) \ !r _s -> do
+         r1 <- (2*) <$> erand48
+
+ ...
-       modifyIORef r (+ rad .* recip (fromIntegral samps))
-       ci <- VM.unsafeRead c i
-       Vec rr rg rb <- readIORef r
-       VM.unsafeWrite c i $ ci + Vec (clamp rr) (clamp rg) (clamp rb) .* 0.25
+       pure (r + rad .* recip (fromIntegral samps))
+       pure (ci + Vec (clamp rr) (clamp rg) (clamp rb) .* 0.25)
+
+ ...
```

## Optimizing smallpt

2020-11-04

### Removing mutation: eliminate IORef

```
Removing mutation: eliminate IORef
...
-   r <- VM.replicate (w * h) 0
-   xi <- newIORef 0
-   flip mapM_ [0..h-1] $ \y -> do
-     writeXi xi y
-     for_ [0..w-1] \x -> do
-       let i = (h-y-1) * w + x
-       for_ [0..1] \sy -> do
-         for_ [0..1] \sx -> do
-           r <- newIORef 0
-           for_ [0..samps-1] \_s -> do
-             r1 <- (2*) <$> erand48 xi
+ img = (`concatMap` [(h-1),(h-2)..0]) $ \y -> runWithErand48 y do
+   for [0..w-1] \x -> do
+     (\pf -> foldlM pf 0 [(sy, sx) | sy <- [0,1], sx <- [0,1]]) \ci (sy, sx) -> do
+       Vec rr rg rb <- (\f -> foldlM f 0 [0..samps-1]) \ !r _s -> do
+         r1 <- (2*) <$> erand48
+
+ ...
-       modifyIORef r (+ rad .* recip (fromIntegral samps))
-       ci <- VM.unsafeRead c i
-       Vec rr rg rb <- readIORef r
-       VM.unsafeWrite c i $ ci + Vec (clamp rr) (clamp rg) (clamp rb) .* 0.25
+       pure (r + rad .* recip (fromIntegral samps))
+       pure (ci + Vec (clamp rr) (clamp rg) (clamp rb) .* 0.25)
+
+ ...
```

All these mutability locations throw in extra RTS code, extra sequencing that blocks the compiler's optimization, and dependency chains. Sometimes we need mutability for performance but in fact one primary optimization technique powering modern compilers is SSA. We almost start there as a functional language don't break it when you don't have a good reason.

## Set everything in smallpt to be strict (1.17x)

```
smallpt :: Int -> Int -> Int -> IO ()
smallpt w h nsamps = do
-   let samps = nsamps `div` 4
-   org = Vec 50 52 295.6
-   dir = norm $ Vec 0 (-0.042612) (-1)
-   cx = Vec (fromIntegral w * 0.5135 / fromIntegral h) 0 0
-   cy = norm (cx `cross` dir) `mulvs` 0.5135
+   let !samps = nsamps `div` 4
+   !org = Vec 50 52 295.6
+   !dir = norm $ Vec 0 (-0.042612) (-1)
+   !cx = Vec (fromIntegral w * 0.5135 / fromIntegral h) 0 0
+   !cy = norm (cx `cross` dir) `mulvs` 0.5135
...
-   r1 <- (2*) `fmap` erand48 xi
-   let dx = if r1<1 then sqrt r1-1 else 1-sqrt(2-r1)
-   r2 <- (2*) `fmap` erand48 xi
-   let dy = if r2<1 then sqrt r2-1 else 1-sqrt(2-r2)
-   d = ...
-   rad <- radiance (Ray (org`addv`(d`mulvs`140)) (norm d)) 0 xi
+   !r1 <- (2*) `fmap` erand48 xi
+   let !dx = if r1<1 then sqrt r1-1 else 1-sqrt(2-r1)
+   !r2 <- (2*) `fmap` erand48 xi
+   let !dy = if r2<1 then sqrt r2-1 else 1-sqrt(2-r2)
+   !d = ...
...
+   pure $! r + rad .* recip (fromIntegral samps)
+   pure $! ci + Vec (clamp rr) (clamp rg) (clamp rb) .* 0.25
```

## Optimizing smallpt

2020-11-04

Set everything in smallpt to be strict (1.17x)

```
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smallpt :: Int -> Int -> Int -> IO ()
smallpt w h nsamps = do
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-   cx = Vec (fromIntegral w * 0.5135 / fromIntegral h) 0 0
-   cy = norm (cx `cross` dir) `mulvs` 0.5135
+   let !samps = nsamps `div` 4
+   !org = Vec 50 52 295.6
+   !dir = norm $ Vec 0 (-0.042612) (-1)
+   !cx = Vec (fromIntegral w * 0.5135 / fromIntegral h) 0 0
+   !cy = norm (cx `cross` dir) `mulvs` 0.5135
...
-   r1 <- (2*) `fmap` erand48 xi
-   let dx = if r1<1 then sqrt r1-1 else 1-sqrt(2-r1)
-   r2 <- (2*) `fmap` erand48 xi
-   let dy = if r2<1 then sqrt r2-1 else 1-sqrt(2-r2)
-   d = ...
-   rad <- radiance (Ray (org`addv`(d`mulvs`140)) (norm d)) 0 xi
+   !r1 <- (2*) `fmap` erand48 xi
+   !dx = if r1<1 then sqrt r1-1 else 1-sqrt(2-r1)
+   !r2 <- (2*) `fmap` erand48 xi
+   !dy = if r2<1 then sqrt r2-1 else 1-sqrt(2-r2)
+   !d = ...
...
+   pure $! r + rad .* recip (fromIntegral samps)
+   pure $! ci + Vec (clamp rr) (clamp rg) (clamp rb) .* 0.25
```

This is not a recommendation, this is a warning. We get a speedup here but it can also regress performance. Some of these bangs are regressions that are hidden. Don't senselessly bang everything in sight.



# Why strictness may be bad

```
let foo = let x = error "ERR" in \y -> y
let foo' = let !x = error "ERR" in \y -> y
```

## Optimizing smallpt

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Why strictness may be bad

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```

# Why strictness may be bad

```
let foo = let x = error "ERR" in \y -> y
let foo' = let !x = error "ERR" in \y -> y

let fooOpt = \y -> y
let foo'Opt = \y -> y  -- ERROR! forcing foo' should give "ERR"
```

## Optimizing smallpt

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let foo = let x = error "ERR" in \y -> y
let foo' = let !x = error "ERR" in \y -> y

let fooOpt = \y -> y
let foo'Opt = \y -> y  -- ERROR! forcing foo' should give "ERR"

let tuple = (42, error "urk")
let bar = let (x, y) = tuple in \z -> x + z
```

## Optimizing smallpt

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Why strictness may be bad

Why strictness may be bad

```
let foo = let x = error "ERR" in \y -> y
let foo' = let !x = error "ERR" in \y -> y
let fooOpt = \y -> y
let foo'Opt = \y -> y  -- ERROR! forcing foo' should give "ERR"

let tuple = (42, error "urk")
let bar = let (x, y) = tuple in \z -> x + z
```

# Why strictness may be bad

```
let foo = let x = error "ERR" in \y -> y
let foo' = let !x = error "ERR" in \y -> y

let fooOpt = \y -> y
let foo'Opt = \y -> y  -- ERROR! forcing foo' should give "ERR"

let tuple = (42, error "urk")
let bar = let (x, y) = tuple in \z -> x + z

let barOpt = \z -> 42 + z
```

## Optimizing smallpt

2020-11-04

Why strictness may be bad

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let foo = let x = error "ERR" in \y -> y
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let fooOpt = \y -> y
let foo'Opt = \y -> y  -- ERROR! forcing foo' should give "ERR"

let tuple = (42, error "urk")
let bar = let (x, y) = tuple in \z -> x + z
let barOpt = \z -> 42 + z
```

# Why strictness may be bad

```
let foo = let x = error "ERR" in \y -> y
let foo' = let !x = error "ERR" in \y -> y

let fooOpt = \y -> y
let foo'Opt = \y -> y  -- ERROR! forcing foo' should give "ERR"

let tuple = (42, error "urk")
let bar = let (x, y) = tuple in \z -> x + z

let barOpt = \z -> 42 + z

let bar' = let (x, !y) = tuple in \z -> x + z
```

## Optimizing smallpt

2020-11-04

Why strictness may be bad

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```
let foo = let x = error "ERR" in \y -> y
let foo' = let !x = error "ERR" in \y -> y
let fooOpt = \y -> y
let foo'Opt = \y -> y  -- ERROR! forcing foo' should give "ERR"

let tuple = (42, error "urk")
let bar = let (x, y) = tuple in \z -> x + z
let barOpt = \z -> 42 + z
let bar' = let (x, !y) = tuple in \z -> x + z
```

# Why strictness may be bad

```
let foo = let x = error "ERR" in \y -> y
let foo' = let !x = error "ERR" in \y -> y

let fooOpt = \y -> y
let foo'Opt = \y -> y -- ERROR! forcing foo' should give "ERR"

let tuple = (42, error "urk")
let bar = let (x, y) = tuple in \z -> x + z

let barOpt = \z -> 42 + z

let bar' = let (x, !y) = tuple in \z -> x + z

let bar'Opt = \z -> 42 + z -- ERROR: forcing bar' should give "urk"
```

## Optimizing smallpt

2020-11-04

Why strictness may be bad

Why strictness may be bad

```
let foo = let x = error "ERR" in \y -> y
let foo' = let !x = error "ERR" in \y -> y
let fooOpt = \y -> y
let foo'Opt = \y -> y -- ERROR! forcing foo' should give "ERR"

let tuple = (42, error "urk")
let bar = let (x, y) = tuple in \z -> x + z
let barOpt = \z -> 42 + z
let bar' = let (x, !y) = tuple in \z -> x + z
let bar'Opt = \z -> 42 + z -- ERROR: forcing bar' should give "urk"
```

## Reduce to only useful strictnesses in smallpt(1.17x)

```
- let !samps = nsamps `div` 4
- !org = Vec 50 52 295.6
- !dir = norm $ Vec 0 (-0.042612) (-1)
- !cx = Vec (fromIntegral w * 0.5135 / fromIntegral h) 0 0
- !cy = norm (cx `cross` dir) .* 0.5135
+ let samps = nsamps `div` 4
+ org = Vec 50 52 295.6
+ dir = norm $ Vec 0 (-0.042612) (-1)
+ cx = Vec (fromIntegral w * 0.5135 / fromIntegral h) 0 0
+ cy = norm (cx `cross` dir) .* 0.5135
...
- !r1 <- (2*) <$> erand48
+ r1 <- (2*) <$> erand48
...
- !r2 <- (2*) <$> erand48
+ r2 <- (2*) <$> erand48
...
- !rad <- radiance (Ray (org+d.*140) (norm d)) 0
+ rad <- radiance (Ray (org+d.*140) (norm d)) 0
...
- pure $! r + rad .* recip (fromIntegral samps)
- pure $! ci + Vec (clamp rr) (clamp rg) (clamp rb) .* 0.25
+ pure (r + rad .* recip (fromIntegral samps))
+ pure (ci + Vec (clamp rr) (clamp rg) (clamp rb) .* 0.25)
```

## Optimizing smallpt

2020-11-04

Reduce to only useful strictnesses in smallpt(1.17x)

```
Reduce to only useful strictnesses in smallpt(1.17x)
- let tsamps = nsamps `div` 4
- !org = Vec 50 52 295.6
- !dir = norm $ Vec 0 (-0.042612) (-1)
- !cx = Vec (fromIntegral w * 0.5135 / fromIntegral h) 0 0
- !cy = norm (cx `cross` dir) .* 0.5135
+ let samps = nsamps `div` 4
+ org = Vec 50 52 295.6
+ dir = norm $ Vec 0 (-0.042612) (-1)
+ cx = Vec (fromIntegral w * 0.5135 / fromIntegral h) 0 0
+ cy = norm (cx `cross` dir) .* 0.5135
...
- !r1 <- (2*) <$> erand48
- !r2 <- (2*) <$> erand48
+ r1 <- (2*) <$> erand48
+ r2 <- (2*) <$> erand48
...
- !rad <- radiance (Ray (org+d.*140) (norm d)) 0
- !rad <- radiance (Ray (org+d.*140) (norm d)) 0
+ rad <- radiance (Ray (org+d.*140) (norm d)) 0
...
- pure $! r + rad .* recip (fromIntegral samps)
- pure $! ci + Vec (clamp rr) (clamp rg) (clamp rb) .* 0.25
+ pure (r + rad .* recip (fromIntegral samps))
+ pure (ci + Vec (clamp rr) (clamp rg) (clamp rb) .* 0.25)
```

Bangs force evaluation. The computation might diverge. Thus the compiler can no longer move the computation around or simplify it. Also if the work doesn't have to get done it still might be. A little thinking about how the variables are used or looking at core allows us to select which ones we bang selectively.

## Use strictness strategically in entire project

```
...
- if det<0 then Nothing else f (b-sdet) (b+sdet)
- where op = p - o
-       eps = 1e-4
-       b = dot op d
-       det = b*b - dot op op + r*r
-       sdet = sqrt det
-       f a s = if a>eps then Just a else if s>eps then Just s else Nothing
+ if det<0
+ then Nothing
+ else
+   let !eps = 1e-4
+       !sdet = sqrt det
+       !a = b-sdet
+       !s = b+sdet
+   in if a>eps then Just a else if s>eps then Just s else Nothing
...
```

## Optimizing smallpt

2020-11-04

### Use strictness strategically in entire project

Use strictness strategically in entire project

```
...
- if det<0 then Nothing else f (b-sdet) (b+sdet)
- where op = p - o
-       eps = 1e-4
-       b = dot op d
-       det = b*b - dot op op + r*r
-       sdet = sqrt det
-       f a s = if a>eps then Just a else if s>eps then Just s else Nothing
+ if det<0
+ then Nothing
+ else
+   let !eps = 1e-4
+       !sdet = sqrt det
+       !a = b-sdet
+       !s = b+sdet
+   in if a>eps then Just a else if s>eps then Just s else Nothing
...
```

Sometimes (point out 'intersect') we have to rearrange the code though when we use bangs. Bangs tell the compiler to make more efficient code, but take away the compiler's options in how to do so. Only take away the compiler's liberties when it's using them poorly. After a little trial and error or reading core how to use bangs in your code will be intuitive.



## Remove Maybe from intersect(s) (1.32x)

| Old: Use Maybe Double to represent (was-hit?:bool, hit-distance: Double)

| New: use (1/0) to represent not (was-hit?)

-intersect :: Ray -> Sphere -> Maybe Double

+intersect :: Ray -> Sphere -> Double

intersect (Ray o d) (Sphere r p \_e \_c \_refl) =

- if det<0 then Nothing else f (b-sdet) (b+sdet)

+ if det<0 then (1/0.0) else f (b-sdet) (b+sdet)

where op = p `subv` o

...

- f a s = if a>eps then Just a else if s>eps then Just s else Nothing

+ f a s = if a>eps then a else if s>eps then s else (1/0.0)

-intersects :: Ray -> (Maybe Double, Sphere)

+intersects :: Ray -> (Double, Sphere)

intersects ray = (k, s)

- where (k,s) = foldl' f (Nothing,undefined) spheres

- f (k',sp) s' = case (k',intersect ray s') of

- (Nothing,Just x) -> (Just x,s')

- (Just y,Just x) | x < y -> (Just x,s')

- \_ -> (k',sp)

+ where (k,s) = foldl' f (1/0.0,undefined) spheres

+ f (k', sp) s' = let !x = intersect ray s' in if x < k' then (x, s') else (k', sp)

radiance :: Ray -> Int -> STRefU s Word64 -> ST s Vec

radiance ray@(Ray o d) depth xi = case intersects ray of

- (Nothing,\_) -> return zerov

- (Just t,Sphere \_r p e c refl) -> do

+ (t,\_) | t == (1/0.0) -> return zerov

+ (t,Sphere \_r p e c refl) -> do

## Optimizing smallpt

2020-11-04

### Remove Maybe from intersect(s) (1.32x)

```
Remove Maybe from intersect(s) (1.32x)
[ Old: Use Maybe Double to represent (was-hit?:bool, hit-distance: Double)
[ New: use (1/0) to represent not (was-hit?)
intersect :: Ray -> Sphere -> Maybe Double
intersect :: Ray -> Sphere -> Double
intersect :: Ray -> Sphere -> Double
intersect (Ray o d) (Sphere r p _e _c _refl) =
- if det<0 then Nothing else f (b-sdet) (b+sdet)
+ if det<0 then (1/0.0) else f (b-sdet) (b+sdet)
where op = p `subv` o
...
- f a s = if a>eps then Just a else if s>eps then Just s else Nothing
+ f a s = if a>eps then a else if s>eps then s else (1/0.0)

intersect :: Ray -> (Maybe Double, Sphere)
intersect :: Ray -> (Double, Sphere)
intersect ray = (k, s)
where (k,s) = foldl' f (Nothing,undefined) spheres
f (k',sp) s' = case (k',intersect ray s') of
- (Nothing,Just x) -> (Just x,s')
- (Just y,Just x) | x < y -> (Just x,s')
- _ -> (k',sp)
+ where (k,s) = foldl' f (1/0.0,undefined) spheres
+ f (k', sp) s' = let !x = intersect ray s' in if x < k' then (x, s') else (k', sp)

radiance :: Ray -> Int -> STRefU s Word64 -> ST s Vec
radiance ray@(Ray o d) depth xi = case intersects ray of
- (Nothing,_) -> return zerov
- (Just t,Sphere _r p e c refl) -> do
+ (t,_) | t == (1/0.0) -> return zerov
+ (t,Sphere _r p e c refl) -> do
```

This is a far more performance critical version of what we saw with 'maximum' vs. 'max'.

Our innermost functions are of critical importance. Here we remove a Maybe which significantly reduces the boxing (which could have been mitigated with a StrictMaybe) and the cases.

Since a Ray that fails to intersect something can be said to intersect at infinity, Double already actually covers the structure at play.

This also reduces allocation.

# Hand unroll the fold in intersects (1.35x)

```
intersects :: Ray -> (Double, Sphere)
-intersects ray = (k, s)
-  where (k,s) = foldl' f (1/0.0,undefined) spheres
+intersects ray =
+  f (... (f (f (intersect ray sphLeft, sphLeft) sphRight) ...)
+  where
    f (k', sp) s' = let !x = intersect ray s' in if x < k' then (x, s') else (k', sp)

-spheres :: [Sphere]
-spheres = let s = Sphere ; z = zeroV ; (.* ) = mulvs ; v = Vec in
-  [ s 1e5 (v (1e5+1) 40.8 81.6)    z (v 0.75 0.25 0.25) DIFF --Left
-  , s 1e5 (v (-1e5+99) 40.8 81.6)  z (v 0.25 0.25 0.75) DIFF --Right
-  ...

+sphLeft, sphRight, ... :: Sphere
+sphLeft  = Sphere 1e5  (Vec (1e5+1) 40.8 81.6)    zeroV (Vec 0.75 0.25 0.25) DIFF
+sphRight = Sphere 1e5  (Vec (-1e5+99) 40.8 81.6)  zeroV (Vec 0.25 0.25 0.75) DIFF
+...
```

# Optimizing smallpt

2020-11-04

## Hand unroll the fold in intersects (1.35x)

```
Hand unroll the fold in intersects (1.35x)

intersects :: Ray -> (Double, Sphere)
intersects ray = (k, s)
  where (k,s) = foldl' f (1/0.0,undefined) spheres
intersects ray =
  f (... (f (f (intersect ray sphLeft, sphLeft) sphRight) ...)
  where
    f (k', sp) s' = let !x = intersect ray s' in if x < k' then (x, s') else (k', sp)

spheres :: [Sphere]
spheres = let s = Sphere ; z = zeroV ; (.* ) = mulvs ; v = Vec in
  [ s 1e5 (v (1e5+1) 40.8 81.6)    z (v 0.75 0.25 0.25) DIFF --Left
  , s 1e5 (v (-1e5+99) 40.8 81.6)  z (v 0.25 0.25 0.75) DIFF --Right
  ...

sphLeft, sphRight, ... :: Sphere
sphLeft  = Sphere 1e5  (Vec (1e5+1) 40.8 81.6)    zeroV (Vec 0.75 0.25 0.25) DIFF
sphRight = Sphere 1e5  (Vec (-1e5+99) 40.8 81.6)  zeroV (Vec 0.25 0.25 0.75) DIFF
...
```

Sadly 'intersects' remains one of our most costly functions even after that improvement so we'll focus on it a bit. A common optimization is loop unrolling. Many compilers do this for us, and there are special versions of it like Duff's Device. Sadly GHC is the one compiler I've used that doesn't. We can though implement every variety of it I know of by hand in Haskell, and this is one of the simplest. I believe we could produce a rule that did it for a given function but I've not actually explored that.

## Custom datatype for intersects parameter passing

Old: Tuple with possibly-unevaluated Double and Sphere

New: Reference to a guaranteed-to-be-evaluated Double and Sphere

```
-intersects :: Ray -> (Double, Sphere)
```

```
+data T = T !Double !Sphere
```

```
+
```

```
+intersects :: Ray -> T
```

```
  intersects ray =  
-   f ( ... f (intersect ray sphLeft, sphLeft) sphRight) ... sphLite  
+   f ( ... f (T (intersect ray sphLeft) sphLeft) sphRight) ... sphLite  
  where  
-   f (k', sp) s' =  
-       let !x = intersect ray s' in if x < k' then (x, s') else (k', sp)  
+   f !(T k' sp) !s' =  
+       let !x = intersect ray s' in if x < k' then T x s' else T k' sp
```

```
radiance :: Ray -> Int -> Erand48 Vec  
radiance ray@(Ray o d) depth = case intersects ray of  
-  (!t,_) | t == 1/0.0 -> return 0  
-  (!t,!Sphere _r p e c refl) -> do  
+  (T t _) | t == 1/0.0 -> return 0  
+  (T t (Sphere _r p e c refl)) -> do  
    let !x = o + d .* t  
        !n = norm $ x - p  
        !nl = if dot n d < 0 then n else negate n
```

## Optimizing smallpt

2020-11-04

### Custom datatype for intersects parameter passing

```
Old: Tuple with possibly-unevaluated Double and Sphere  
New: Reference to a guaranteed-to-be-evaluated Double and Sphere  
-intersects :: Ray -> (Double, Sphere)  
-data T = T !Double !Sphere  
+  
+intersects :: Ray -> T  
+intersects ray =  
+  f (...) f (intersect ray sphLeft, sphLeft) sphRight) ... sphLite  
+  f (...) f (T (intersect ray sphLeft) sphLeft) sphRight) ... sphLite  
+  where  
+  - f (k', sp) s' =  
+  -   let !x = intersect ray s' in if x < k' then (x, s') else (k', sp)  
+  + f !(T k' sp) !s' =  
+  +   let !x = intersect ray s' in if x < k' then T x s' else T k' sp  
  
radiance :: Ray -> Int -> Erand48 Vec  
radiance ray@(Ray o d) depth = case intersects ray of  
-  (!t,_) | t == 1/0.0 -> return 0  
-  (!t,!Sphere _r p e c refl) -> do  
+  (T t _) | t == 1/0.0 -> return 0  
+  (T t (Sphere _r p e c refl)) -> do  
    let !x = o + d .* t  
        !n = norm $ x - p  
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```

Still trying to focus on the performance of 'intersects' we can optimize how it passes data to itself and its caller. We want the data to be strict so the compiler knows the tuple's members will always be evaluated, but we want to avoid copying. A normal tuple lacks strictness information. An unboxed tuple sadly forces copying because of its unpacked nature. Sphere is quite large, and will be expensive to copy. What we need is a strict tuple, and we create one for just this usage. This exists in libraries of course, but we wanted to illustrate it.

# Optimize file writing

```
build-depends:
    base >= 4.12 && < 4.15
+    , bytestring ^>= 0.11

-toInt :: Double -> Int
-toInt x = floor $ clamp x ** recip 2.2 * 255 + 0.5
+toInt :: Double -> BB.Builder -- O(1) concatenation
+toInt x = BB.intDec (floor (clamp x ** recip 2.2 * 255 + 0.5)) <> BB.char8 ' '
...
withFile "image.ppm" WriteMode $ \hdl -> do
-    hPrintf hdl "P3\n%d %d\n%d\n" w h (255::Int)
-    for_ img \(Vec r g b) -> do
-        hPrintf hdl "%d %d %d " (toInt r) (toInt g) (toInt b)
+    BB.hPutBuilder hdl $
+        BB.string8 "P3\n" <> -- efficient builders for ASCII
+        BB.intDec w <> BB.char8 ' ' <> BB.intDec h <> BB.char8 '\n' <>
+        BB.intDec 255 <> BB.char8 '\n' <>
+        (mconcat $ fmap \(Vec r g b) -> toInt r <> toInt g <> toInt b) img)
```

## Optimizing smallpt

2020-11-04

### Optimize file writing

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Optimize file writing

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+        BB.intDec w <> BB.char8 ' ' <> BB.intDec h <> BB.char8 '\n' <>
+        BB.intDec 255 <> BB.char8 '\n' <>
+        (mconcat $ fmap \(Vec r g b) -> toInt r <> toInt g <> toInt b) img)
```

We're getting quite a bit faster now, but there's still some small pieces of low lying fruit. Strings are fairly inefficient and the conversion functions to them we use aren't particularly efficient. 'bytestring' has some efficient writing code, so we just convert to that for a modest gain.

Use LLVM backend (1.87x)

```
+package smallpt-opt
+  ghc-options: -fllvm
```

## Optimizing smallpt

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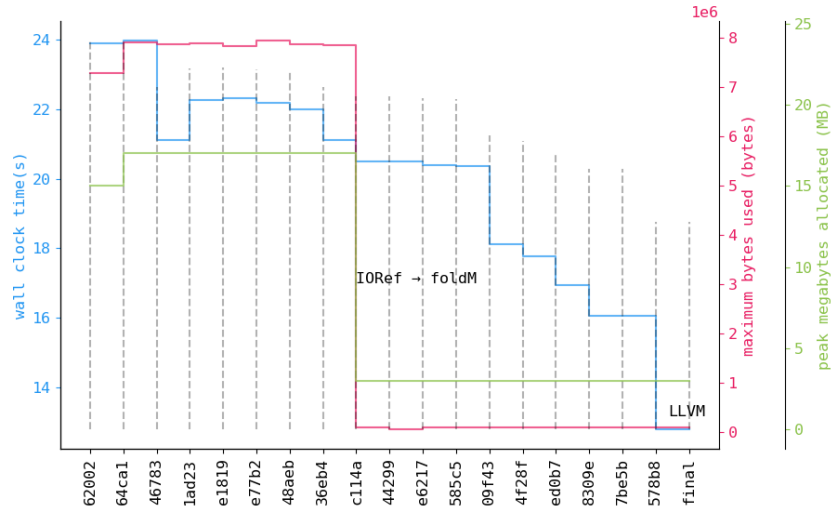
└ Use LLVM backend (1.87x)

Use LLVM backend (1.87x)

```
+package smaller-opt
+  ghc-options: -fllvm
```

Finally, this particular code is quite numeric heavy. There are optimizations for numeric heavy code we're missing in GHC. LLVM has an extensive library of laws to optimize low level numeric ops. LLVM is too low-level to understand haskell as haskell. It makes decisions with the tacit assumption that the assembly came from a C-like language, which is often to the detriment of a Haskell-like language. In this case, as the code is "fortran-like", LLVM wins.

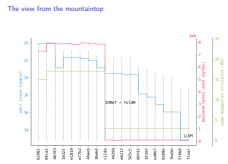
# The view from the mountaintop



## Optimizing smallpt

2020-11-04

The view from the mountaintop



# Takeaways

- ▶ The unrolling in 'intersects' is ugly.
- ▶ (We feel) the maintainability of this code hasn't been significantly harmed.
- ▶ We're faster than clang++ and within 6% of g++
- ▶ Haven't exhausted the optimization opportunities.
- ▶ GHC could learn to do several of these optimizations for us.
- ▶ Others are just good Haskell style.
- ▶ Clean Haskell is often performant Haskell.

## Optimizing smallpt

2020-11-04

### Takeaways

The unrolling in 'intersects' is ugly but other than that optimization it is our opinion that the maintainability of this code hasn't been significantly harmed by improving its performance. We're faster than clang++ and within 6% of g++ We haven't exhausted the optimization opportunities. I believe GHC could learn to do several of these optimizations for us. Many of the others are just good Haskell style. Clean Haskell is often performant Haskell.

Takeaways

- ▶ The unrolling in 'intersects' is ugly.
- ▶ (We feel) the maintainability of this code hasn't been significantly harmed.
- ▶ We're faster than clang++ and within 6% of g++
- ▶ Haven't exhausted the optimization opportunities.
- ▶ GHC could learn to do several of these optimizations for us.
- ▶ Others are just good Haskell style.
- ▶ Clean Haskell is often performant Haskell.