

High Performance Iterated Function Systems

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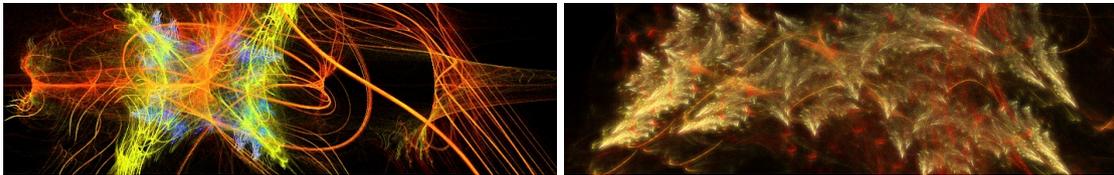


Figure 1: A crop from a Fractal Flame image rendered using $2 \cdot 10^6$ points at a resolution of 1280×800 with 40 Hz on a GTX280.

This article presents an implementation of the Fractal Flames algorithm. It uses CUDA and OpenGL to take advantage of the computational power of modern graphics cards. GPUs use a SIMT (Single Instruction Multiple Thread) architecture. It is needed to design programs in a way which avoids divergent branching. The Fractal Flames algorithm involves random function selection which needs to be calculated in each thread. The algorithm thus causes heavy branch divergence which leads to $O(n)$ complexity in the number of functions. Current implementations suffer severely from this problem and address it by restricting the amount of functions to a value with acceptable computational overhead.

The implementation presented in this paper changes the algorithm in a way that leads to $O(1)$ complexity and therefore allows an arbitrary amount of active functions. This is done by applying a presorting algorithm that removes the random function selection and thus eliminates branch divergence completely.

1 Problem Statement and Mathematical Background

Fractal Flames is an algorithm to create fractal images based on Iterated Function Systems (IFS) with a finite set of functions. The algorithm uses the *Chaos Game* [Bar88], which is an iteration scheme that picks one random function for each data point and iteration, evaluates it, and continues with the next iteration. This scheme is visualized in Figure 2. The larger

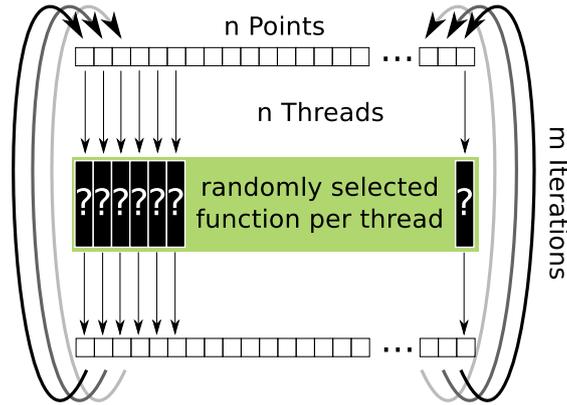


Figure 2: Naive parallelizing approach of the Chaos Game. The parallelization is realized by assigning a thread to each data point. Every thread chooses a random function and replaces the point by the computed value. This process is repeated m times.

the number of different functions, the more interesting the images that can be rendered. It is therefore a goal to allow for as many functions as possible.

Selecting a random function per sample results in a different branch per thread, if the algorithm is implemented on graphics hardware. One needs to consider that these devices execute the same instruction in lockstep within a *warp* of 32 threads. That is, the multiprocessor can execute a single branch at a time, and therefore diverging branches are executed in serial order.

The random function selection causes very heavy branch divergence as every thread needs to evaluate a different function in the worst case. This results in linear runtime complexity in the number of functions which severely restricts the number of functions that can be used in practice, due to the computational overhead involved.

1.1 Iterated Function Systems

As described in [Bar88], an IFS consists of a finite set of affine contractive functions

$$F = \{f_i : X \rightarrow X \mid i = 1, 2, \dots, N\}, N \in \mathbb{N}. \quad (1)$$

The set $F_s = \{f_1^s, f_2^s, f_3^s\}$ for example forms a Sierpinski triangle with

$$f_1^s(p) = \frac{p}{2}, \quad f_2^s(p) = \frac{1}{2} \left(p + \begin{pmatrix} 0 \\ 1 \end{pmatrix} \right), \quad f_3^s(p) = \frac{1}{2} \left(p + \begin{pmatrix} 1 \\ 0 \end{pmatrix} \right), \quad (2)$$

where p is a vector.

The associated set S is the fixed point of Hutchinson's recursive set equation

$$S = \bigcup_{i=1}^N f_i(S). \quad (3)$$

It is not possible to directly evaluate the set, as the Hutchinson equation (Equation 3) describes an infinite recursion. An approximate approach is the Chaos Game [Bar88] which solves the Hutchinson equation by a Monte Carlo method. Figure 3 shows it in a basic version. The Chaos Game starts by selecting a random point off the bi-unit square $p = (x, y)$ with $|x| \leq 1, |y| \leq 1$ and starts its iterating phase by choosing a random function f_i each iteration and evaluates $p := f_i(p)$. With increasing number of calculated iterations, p converges closer to the set [Sil98]. After a sufficient number of steps, the point can be plotted in every iteration.

Iterated Function Systems have numerous applications, as outlined in [FF90, Chapter 9–18], such as graphs of functions, dynamical systems and brownian motion. Furthermore, this approach could be used in multilayer material simulations in ray tracing. In this article we concentrate on IFS for the simulation of Fractal Flames.

We chose to start with a random point and do random function selection, even though randomness is not a requirement of the algorithm. The initial point is actually arbitrary, it will still converge to the set, given that the last iterations contain all possible combinations of functions. As we aim for real time performance, we can only evaluate a very limited number of iterations, but on a lot of points due to parallelization. Should the resulting fractal be too complicated to achieve sub-pixel convergence, the random initialization still creates a uniform appearance. This is, especially over the course of some frames, more visually pleasing than deterministic artifacts.

1.2 Fractal Flames

Fractal Flames, as described in [DR04] extend the IFS algorithm by allowing a larger class of functions that can be in the set F . The only restriction that is imposed on the functions f_i is the contraction on average. These functions can be described by

$$f_i(x, y) = P_i \left(\sum_j v_{ij} V_j (a_i x + b_i y + c_i, d_i x + e_i y + f_i) \right) \quad (4)$$

$p =$ a random point in $[-1, 1]^2$

```

iterate  $m$  times {
   $i :=$  random integer from 0 to  $N - 1$  inclusive
   $p := f_i(p)$ 
  if  $iteration > 17$ 
    plot( $p$ )
}

```

Figure 3: The Chaos Game Monte Carlo algorithm. It chooses a random point and starts its iteration phase. Every iteration, a random function f_i is selected and evaluated for p which is then assigned to p .

with $P_i(x, y) = (\alpha_i x + \beta_i y + \gamma_i, \delta_i x + \epsilon_i y + \zeta_i)$, where $a_i \dots g_i$ and $\alpha_i \dots \zeta_i$ express affine transformations on 2D points, while V_j , so-called *variations* apply non-linear transformations, which are scaled by the factors v_{ij} . Typically, each function has its own set of up to 20 variations V_j . A few variation functions can be found in Figure 4. An extensive collection of those variations can be found in [DR04].

$$\begin{aligned} \psi & : \text{ random number } \in [0, 1], \quad \Omega : \text{ random number } \in [0, \pi] \\ r & = \sqrt{x^2 + y^2}, \quad \theta = \arctan(x/y) \\ V_0(x, y) & = (x, y) \\ V_1(x, y) & = (\sin x, \sin y) \\ V_2(x, y) & = \frac{1}{r^2}(x, y) \\ V_3(x, y) & = (x \sin(r^2) - y \cos(r^2), x \cos(r^2) + y \sin(r^2)) \\ V_{13}(x, y) & = \sqrt{r} \cdot (\cos(\theta/2 + \Omega), \sin(\theta/2 + \Omega)) \\ V_{18}(x, y) & = e^{x-1} \cdot (\cos(\pi y), \sin(\pi y)) \\ V_{19}(x, y) & = r^{\sin \theta} \cdot (\cos \theta, \sin \theta) \end{aligned}$$

Figure 4: A few selected variation functions. Ω and ψ are new random numbers in each evaluation of the variation function.

Every function is assigned a weight w_i which controls the probability that f_i is chosen in a Chaos Game iteration. This parameter controls the influence of a function in the computed image. Furthermore, a color $c_i \in [0, 1]$ is assigned. Every point has a third component which holds the current color c which is updated by $c := (c + c_i)/2$ in each iteration and is finally mapped into the output color space.

The computed points are visualized by creating a colored histogram. Since the computed histogram has a very high dynamic range, a tone mapping operator is applied.

2 Core Technology

In order to remove branch divergence, we replace the randomness of the function selection by randomized data access. This way, instructions can be optimally and statically assigned to threads.

Warps are assigned to a fixed function and every thread randomly selects a data point in each iteration. This selection is realized by a random bijective mapping between the data and the thread indices. A fixed set of precomputed permutations is used as they don't depend on dynamic data and may be cyclic as it doesn't matter if images repeat after a few rendered frames.

Every thread calculates its assigned function and indirectly accesses the data array by its permuted index. It then evaluates the function and writes the data back. A new permutation

is picked in each iteration.

Figure 5 shows the iteration scheme.

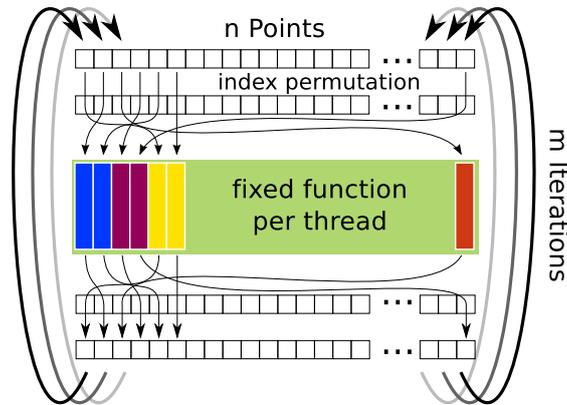


Figure 5: Optimized parallel algorithm. Instead of indexing directly into the data array as in Figure 2, the data is shuffled randomly in every iteration. This allows to statically assign threads to functions and thereby remove the branch divergence.

3 Implementation

The optimized algorithm as well as the divergent approach have been implemented to benchmark both of them. The implementation uses CUDA to be able to tightly control the thread execution which would not have been possible with traditional shading languages.

All variation functions have been implemented in a large switch statement. A struct containing variation indices and the scalar factor is stored in the constant memory. To evaluate a function f_i , a loop over all those indices is performed which is used to index in the switch statement. The results are summed up according to Equation (4).

3.1 The Three Phases

The optimized Chaos Game algorithm requires synchronization across all threads and blocks after each iteration, because the data gets shuffled across all threads. Due to CUDA's lack of such a synchronization instruction, the Chaos Game had to be split into three kernels to achieve synchronization by multiple kernel executions (see Figures 7 and 6):

The initialization kernel performs one iteration of the Chaos Game on the randomly permuted Hammersley point set. The warm-up kernel performs one Chaos Game iteration and is called multiple times until the output data points converged to the fractal. The last kernel generates a larger set of points by randomly transforming each of the previously generated points 64 times, producing 64 independent samples for each input sample. The generated points are recorded in a vertex buffer object (VBO), which is then passed to the rendering

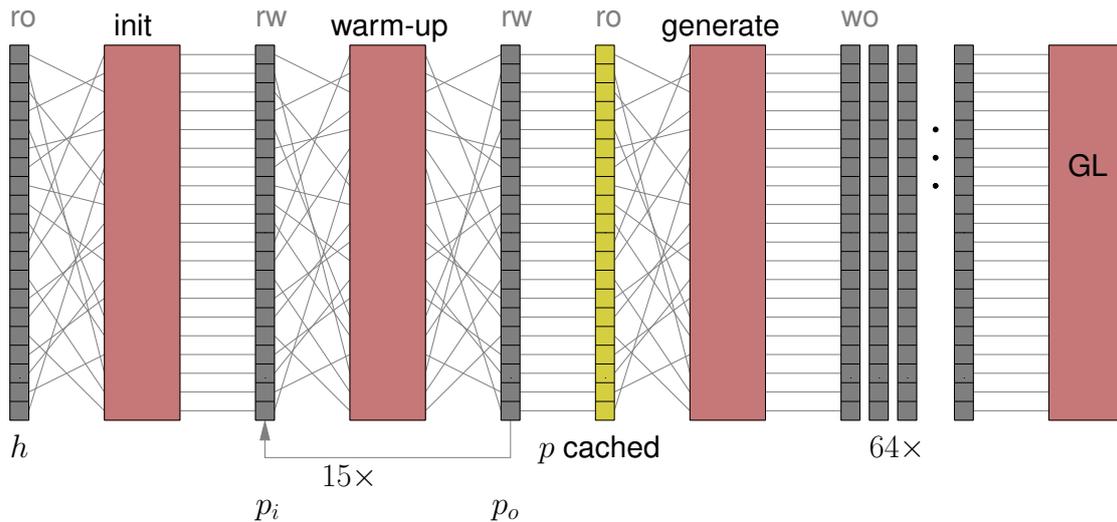


Figure 6: Illustration of the pseudo code listing in Figure 7, the three left red boxes are the three kernels, the last represents rendering in OpenGL. Memory access patterns are indicated by the gray lines. On the left, h indicates the Hammersley point set as the input.

stage. From those kernels, the point generation kernel takes by far the most computation time. Those run times on a GTX280 are $34.80\mu s$ init, $43.12\mu s$ warm-up, and $1660.59\mu s$ generate points. The init kernel occupies 1.5% of the total Chaos Game runtime, the warm-up kernel is called 15 times and thus takes 27.6%, whereas the point generation kernel takes 70.9%.

3.2 Memory Access Patterns

The point generation kernel needs to select a point from the point buffer multiple times, calculate one iteration and write the result into the VBO. As the input point array does not change, texture caching can be used. The effect of this optimization is 6.9% speed improvement for the 9500GT and 18% for the GTX280. The access pattern is pretty bad for the texture caching mechanisms of a GPU as it completely lacks locality. As the memory consumption of the samples is approximately 200kb, the cache of a GPU seems to be big enough to hold a large amount of samples and therefore this method yields a major performance improvement. Benchmarks have been conducted to measure the effect of the number of samples on the texture caching. As shown in Figure 8, the number of threads – and therefore the number of samples in the texture cache – doesn't have a huge impact: The runtime stays constant up to 2^{12} threads and then increases linearly with the number of threads; the cached version is constantly faster.

Note that only the warm-up kernel has a completely random memory access pattern during read and write. It needs to randomly access the point array to read a point for the iteration and write out its result. To avoid race conditions on write accesses, either data has to be written back to the position the point was read from, or else a second point array is

```

f : fixed function for each thread
i : thread index
h := Hammersley point set in the bi-unit square
p : the point buffer
 $\phi$  := random permutation

// phase one: initialization
p[i] := f(h[ $\phi$ (i)])

// phase two: warm-up
iterate 15 times {
     $\phi$  := next random permutation
    p[ $\phi$ (i)] := f(p[ $\phi$ (i)])
}

// phase three: generate points
iterate 64 times {
     $\phi$  := next random permutation
    plot(f(p[ $\phi$ (i)]))
}

```

Figure 7: The three phases of the GPU version of the Chaos Game. First, p is initialized by transforming the Hammersley points by one function evaluation. Second, the warm-up phase assures that the points p are sufficiently close to the fractal. Finally, the point set is transformed 64 times more. In this last phase, the transformed points are displayed, but not written back to p .

needed. The former approach involves random reads and writes, whereas the latter one has higher memory use. The initialization kernel needs to randomly select a start point, but it can write the iterated points in coalesced manner. Due to its low runtime, there is no need to optimize it further.

3.3 Rendering

The generated points are rendered as a tone mapped histogram using OpenGL. This step consumes approximately 70% of the total time, depending on the particular view.

The 1D color component needs to be mapped into the RGBA representation. This is done by a 1D texture look-up in a shader. To create the colored histogram, the points are rendered using the OpenGL point primitive and using an additive blending mode. This yields brighter spots where multiple points are drawn in the same pixel. To cope with the large range of values, a floating point buffer is employed. The resulting histogram is then tone mapped and gamma corrected because the range between dark and bright pixel values is very high. For

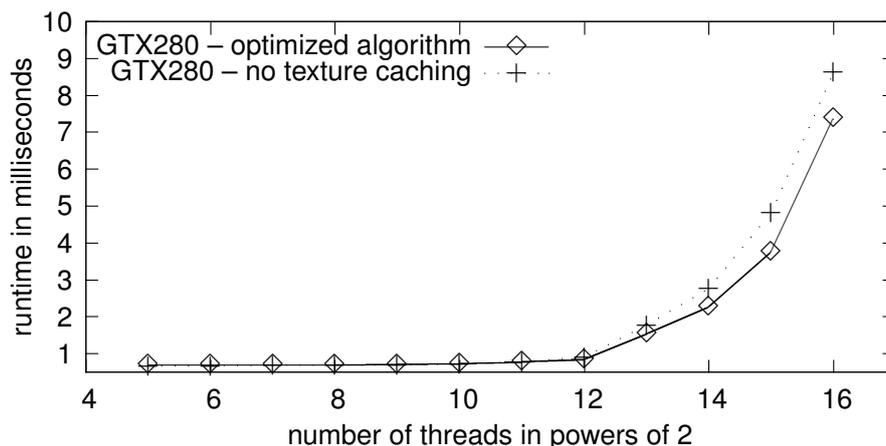


Figure 8: Runtime on the GTX280 in dependence of number of running threads. Each thread generates 64 samples. The runtime stays almost constant to 2^{12} threads and subsequently increases linearly with the number of threads.

sake of simplicity and efficiency, we use $L' = L/(L + 1)$ to calculate the new pixel lightness L' from the input lightness L , but more sophisticated operators are possible [DR04].

Monte Carlo generated images are typically quite noisy. This is especially true for the real-time implementation of Fractal Flames as only a limited number of samples can be calculated in the given time frame. Using temporal anti aliasing smoothes out the noise by blending between the actual frame and the previously rendered ones. This is implemented by using Frame Buffer Objects and increases the image quality significantly.

3.4 Static Data

Some static data is needed during runtime which is calculated by the host during program startup. A set of Hammersley Points [Lem09] is used as the starting point pattern to ensure a stratified distribution, which increases the quality of the generated pictures. The presorting algorithm needs permutations which are generated by creating multiple arrays containing the array index and a random number calculated by the Mersenne Twister [MN98]. These arrays are then sorted by the random numbers, which can be discarded after that. The resulting set of permutations is uploaded onto the device. As some of the implemented functions need random numbers, also a set of random numbers is created and uploaded onto the device. All of this generated data is never changed during runtime.

In the runtime phase, it is necessary to specify various parameters which characterize the fractal. Every function needs to be evaluated in $p(f_i) \cdot num_threads$ threads, where $p(f_i)$ is the user specified probability that function f_i is chosen in the Chaos Game. As the number of functions is small compared to the number of threads, it doesn't matter if this mapping is done in warp size granularity instead of thread granularity. Each thread is mapped to a function by an index array which is compressed into a prefix sum. A thread

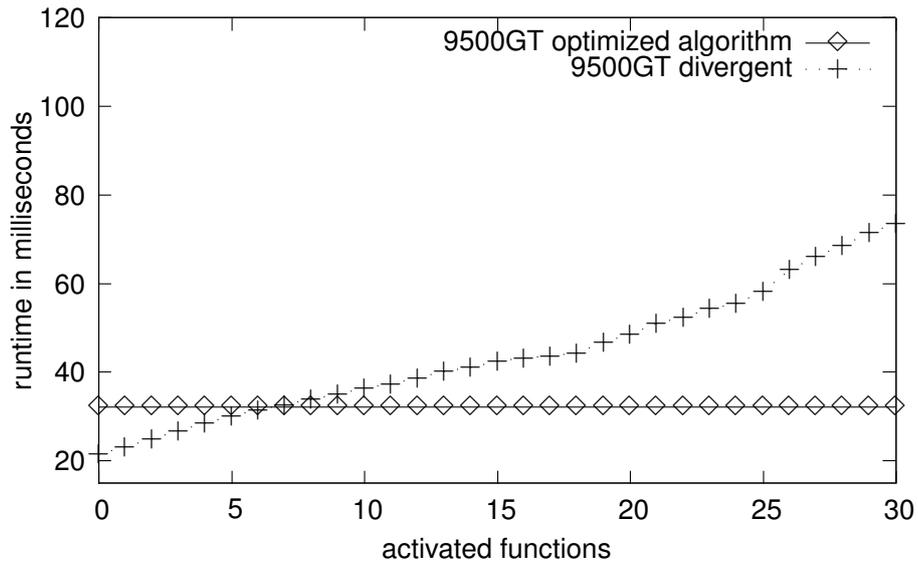


Figure 9: Runtime comparison between divergent and optimized Chaos Game on a 9500GT in dependence on the number of activated functions. The divergent implementation shows linearly increasing runtime in the number of activated functions while the optimized algorithm shows constant runtime. The optimized version beats the divergent solution when eight or more functions are activated.

can find out which function it has to evaluate by performing a binary search on this array. This search does not introduce a performance penalty in our experiments, but reduces the memory requirements significantly. The parameters and the prefix sum are accumulated in a struct that is uploaded into the constant memory each frame.

4 Final Evaluation

The naive parallel implementation degenerates to an $O(n)$ algorithm in the number of functions on graphics hardware due to the branch divergence issues. The proposed algorithm turns the function evaluation in the naive Chaos Game algorithm to an $O(1)$ solution. This is shown in Figure 9 and Figure 10 where the optimized algorithm is benchmarked against the naive parallel implementation on a 9500GT and a 280GTX, respectively. In the benchmark, 20 randomly selected variation functions V_j were used in each f_i . The evaluation time of the optimized algorithm stays constant, whereas the naive solution increasingly suffers from branch divergence issues. Employing the CUDA profiler, the branch divergence can be measured. The amount of branch divergence with 20 activated functions is at 23.54% for the divergent approach, whereas the optimized algorithm only shows branch divergence inside the function evaluation (0.05%).

The optimization doesn't come without cost though. The algorithm basically trades branch

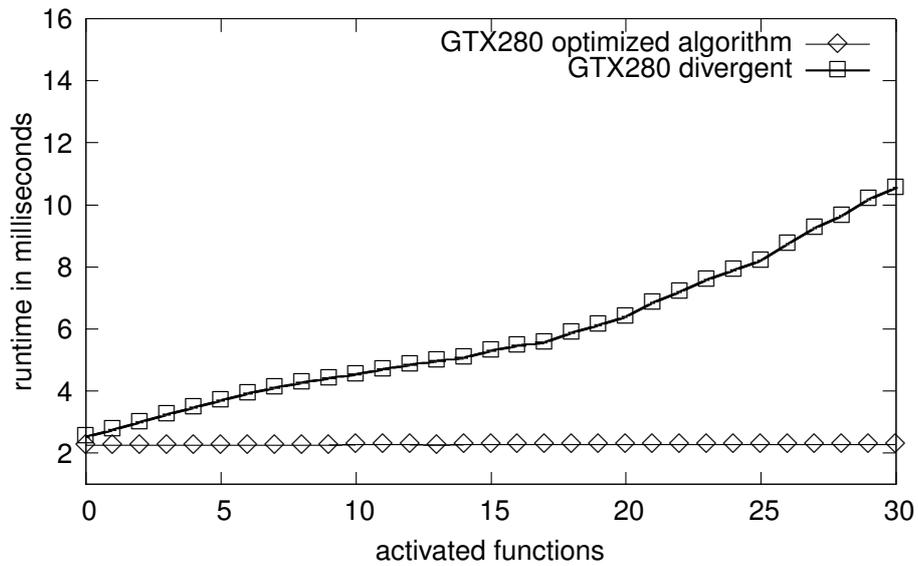


Figure 10: Runtime comparison between divergent and optimized Chaos Game on a GTX280 in dependence on the number of activated functions. The performance characteristics are the same as with the 9500GT shown in Figure 9, with a smoother increase in the divergent graph. The optimized version is always faster.

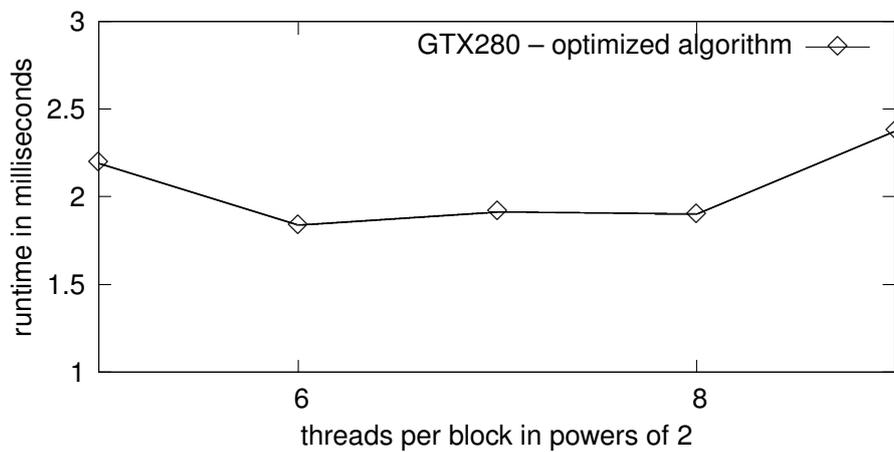


Figure 11: Runtime changing with increasing block size and constant number of threads. The fastest configuration is a block size of 2^6 .

divergence for completely random memory access patterns. As the benchmark shows, GPUs handle this surprisingly well. Different GPUs have been benchmarked to show that this behavior isn't restricted to high end GPUs (the speed impact of permuted read and write access versus completely linear operations was 3.1% for the 9500GT and 5.3% for the 280GTX).

In Figure 11, it is shown that the Chaos Game runtime is not monotonically decreasing with increasing block size. The best runtime is achieved with 64 threads per block. It is not entirely clear from where this behavior arises, but it may be that the scheduler has a higher flexibility with lower block sizes.

In our implementations the occupancy as reported by the CUDA profiler is 0.25. This is due to high register usage of the function evaluation code. Enforcing a lower register usage at compile time increases the occupancy to 1.0 but reduces the performance by 51.3%.

We additionally implemented a CPU version to verify the benefits of using graphics hardware. Benchmarks have shown that a GTX285 is able to perform the Chaos Game iterations without the rendering about 60 times faster than a single-core Opteron CPU clocked at 2.8GHz (240Hz vs. 3.7Hz).

5 Conclusion

A presorting approach has been presented that completely eliminates branch divergence in Monte Carlo algorithms that use random function selection and evaluation. Furthermore it was shown that the approach is sufficiently fast for real-time rendering at high frame rates, while it is necessary to keep an eye on the imposed performance penalty due to high memory bandwidth costs.

The optimized approach allows us to change all parameters in real-time and observe their effects, it thus can effectively help in getting an understanding of Iterated Function Systems. This also makes it possible to use the Fractal Flame algorithm in live performances by changing the parameters in real-time, conducting interactive animations.

References

- [Bar88] M. Barnsley. *Fractals Everywhere: The First Course in Deterministic Fractal Geometry*. Academic Press, 1988.
- [DR04] S. Draves and E. Reckase. The fractal flame algorithm, 2004.
- [FF90] K. Falconer and K.J. Falconer. *Fractal geometry: mathematical foundations and applications*. Wiley New York, 1990.
- [Lem09] C. Lemieux. *Monte Carlo and Quasi-Monte Carlo Sampling*. Springer Verlag, 2009.

- [MN98] M. Matsumoto and T. Nishimura. Mersenne twister: a 623-dimensionally equidistributed uniform pseudo-random number generator. *ACM Transactions on Modeling and Computer Simulation (TOMACS)*, 8(1):3–30, 1998.
- [Sil98] B. Silverman. *Density estimation for statistics and data analysis*. Chapman & Hall/CRC, 1998.