number of pixels written to the frame buffer. The number of rendering passes of the depth-peeling algorithm is the number of layers plus one, which typically is significantly smaller than the number of slices.

When rendering order-independent transparent objects, all the layer images are blended in depth order. In contrast, for voxelization, we want the layers to be separated, which is similar to a layered depth image (Shade et al. 1998). The pixels of the layered images are the attributes of the corresponding voxels. The first attribute is the 3D position, and the other attributes depend on the application. Assume that the maximum size along any of the major axes of the object being voxelized is D. We allocate an off-screen buffer with width and height equal to the maximum number of layers times D and the number of attributes times D, respectively. Then, between rendering passes, we translate the viewport so that the layer images do not overlap but are tiled as tightly as possible. We apply the peeling process three times; each time, the image plane is orthogonal to one of the major axes. That is, we perform the peeling from three orthogonal views to avoid missing voxels. As a result, some of the voxels may be rendered more than once. However, the replication does not affect the accuracy. The images containing the voxel attributes are then copied to a vertex array allocated in video memory (using OpenGL extensions such as `ARB_pixel_buffer_object` and `ARB_vertex_buffer_object` [NVIDIA 2004]). Note that different types of voxel attributes are copied to different locations inside the vertex array. We may avoid the copying if either render-to-vertex-array or vertex texture fetch is available. Figure 47-6 illustrates the process in 2D. The line segments are the boundary lines (surfaces in 3D). The small red, green, and blue boxes represent the voxels generated when the boundary lines are projected onto the layered off-screen buffers. Note that the red boundary will only result in two voxels if it is only rendered into layer X.

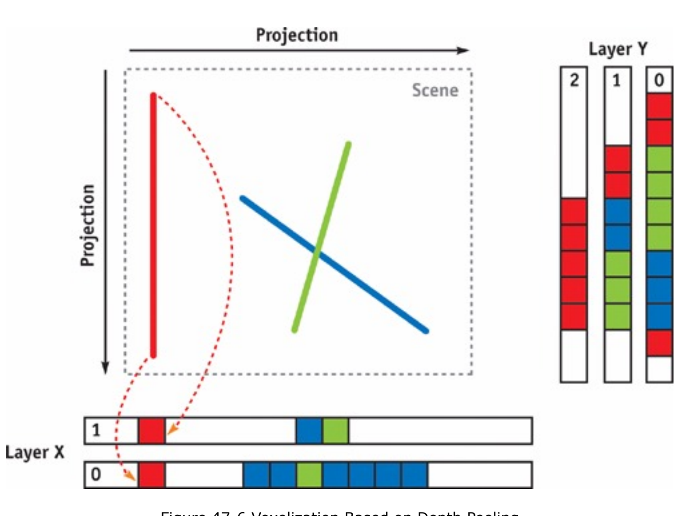

Figure 47-6 Voxelization Based on Depth Peeling

The vertex array is essentially an array of voxel positions and other attributes, which can generate all the boundary voxels for further processing. We use the vertex array because current GPU fragment programs cannot scatter (that is, the output location of a fragment is fixed). Therefore, we must rely on the transformation capability of vertex programs. We convert the vertex array into a flat volume by rendering each vertex as a point of size 1, so that each voxel has a footprint of 1 pixel. All the vertices of the array pass through a vertex program that translates each voxel properly according to its *z* coordinate, using equations similar to Equation 1. The frame buffers for the depth peeling are initialized with large numbers. If a pixel is not covered by any boundary voxels, then the corresponding vertex created from the pixel falls far away from the view frustum and is clipped.

47.4.2 Periodic Boundaries

For periodic boundaries, outgoing packet distributions wrap around and reenter the lattice from the opposite side. In practice, periodic boundaries are actually computed during the propagation of the packet distributions. If a periodic boundary face is orthogonal to the *x* or *y* axis, we call it an *in-slice periodic boundary face*, because a distribution on the face is copied to the opposite side of the lattice but stays inside the same *xy* slice. For in-slice periodic boundaries, we simply apply a modulo operation to the texture coordinates by the width or the height of each slice. If a periodic boundary face is perpendicular to the *z* axis, we call it an *out-slice periodic boundary face*, for which we need to copy distribution textures of one slice to another.

A naive implementation of the in-slice periodic boundary condition is to apply the modulo operation to the texture coordinates of all the distribution texels. However, this can be very slow. Therefore, one optimization we use first propagates without the periodic boundary condition. Then we draw strips of single-texel width that only cover the boundary and apply the modulo operation. (See Chapter 34 of this book, "GPU Flow-Control Idioms," for more information on this technique.) In this way, the cost for computing is negligible, because the periodic boundary nodes account for only a small percentage of the lattice.

47.4.3 Outflow Boundaries

For outflow conditions, some packet distributions of the boundary nodes propagate outside of the lattice and can simply be discarded. However, these nodes also expect some packet distributions to propagate inward from fictitious nodes just outside the boundary. One solution is to copy the distribution of the internal nodes to those boundary nodes, according to Equations 2, 3, and 4, in which $f(B)(i, j, k)$ is the distribution of a boundary node in the velocity direction of (i, j, k); $f(I)(i, j, k)$ is the distribution of the corresponding internal node; and $i, j, k \in \{-1, 0, 1\}$. For example, boundary nodes on slice 0 are copied from internal nodes two slices away, which are on slice 2.

For an outflow face orthogonal to the *x* axis:

Equation 2

$$f(B)_{(i,j,k)} = f(I)_{(-i,j,k)}.$$

For an outflow face orthogonal to the *y* axis:

Equation 3

$$f(B)_{(i,j,k)} = f(I)_{(i,-j,k)}.$$

For an outflow face orthogonal to the *z* axis:

Equation 4

$$f(B)_{(i,j,k)} = f(I)_{(i,j,-k)}.$$

Similar to periodic boundaries, an outflow boundary condition is applied by drawing single-texel-wide stripes. Note that when packing the distributions, we guarantee that $f_{(i, j, k)}$ and $f_{(-i, j, k)}$ of the same node are in the same texture, as well as the pairs $f_{(i, j, k)}$ and $f_{(i, -j, k)}$, $f_{(i, j, k)}$ and $f_{(i, j, -k)}$. You can easily verify these with Table 47-1. Therefore, each boundary distribution texture copies from only one distribution texture. To flip the distributions around the major axes, we use the swizzling operator to rearrange the order of the color elements.

47.4.4 Obstacle Boundaries

For obstacle boundaries, we adopt the improved bounce-back rule of Mei et al. 2000, which can handle deformable and moving boundaries. As indicated in Figure 47-7, all packet distributions with corresponding links that are intersected by a boundary are updated using the curved boundary equations rather than the standard LBM equations. For example, the link between the two nodes in Figure 47-7 is intersected by a boundary, hence the boundary rules are applied to the packet distributions marked in blue, f_1 and f_2. Again, we don't need to worry about the details of these equations; interested readers will find these equations in Mei et al. 2000. We just need to know that the boundary equations require two values in addition to the values of f_1 and f_2 at the previous time step. The first value is the exact intersection location, which is described by the ratio D, where D = $t_1/(t_1 + t_2)$ for f_1, and D = $t_2/(t_1 + t_2)$ for f_2. The second value is the moving speed at the intersection, $\mathbf{u}_w$, which is useful for moving and deformable boundaries. Note that for each boundary link, the boundary condition affects two distributions, one on each side of the boundary. The two distributions are in opposite directions, but they are colinear with the link. We refer to them as *boundary distributions*.

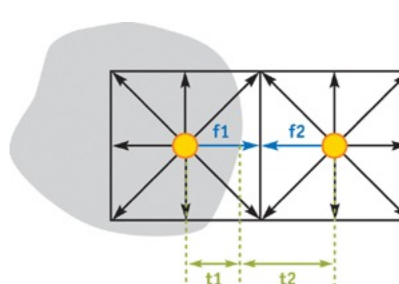

Figure 47-7 Curved-Wall Boundary in LBM Lattice

To generate the boundary information, we first create a voxelization of the boundaries using the method described in the previous section. Besides the position of the voxels, we also need the wall velocity $\mathbf{u}_w$, as well as the coefficients of polygon plane equations, which will be used to compute D. That is, we need three attributes in total. To preserve accuracy, we treat these attributes as texture coordinates when rendering the vertex array in the next step.

In practice, we don't explicitly generate the flat volume of the boundary voxels. Rather, we combine the generation with the computation of the boundary conditions, by rendering the boundary vertex array directly into the off-screen buffer containing the propagated new distributions and then applying the fragment program for complex boundary conditions. Note that in most cases only one node of each boundary link is covered by a voxel from the generated voxel array. However, we need each boundary node to receive a fragment so that the boundary distributions are updated. Therefore, for each packed distribution texture, we render the voxel array three times. In the first pass, they are rendered normally, covering those voxels in the generated voxel array. In the second pass, we first set the color mask so that only the R and G channels can be modified. Then we apply a translation to all the voxels using a vertex program. The translated offset is computed according to the following rule:

$$\cdot\ \mathbf{e}_{qi} \quad : \text{if } flag_1 \times flag_2 > 0;$$

$$\cdot -\mathbf{e}_{qi} \quad : \text{if } flag_1 \times flag_2 < 0;$$

$$\cdot\ 0 \quad : \text{if } flag_1 \times flag_2 = 0;$$

where $flag_1 = (pos_x, pos_y, pos_z, 1) \cdot (A, B, C, D)$, $flag_2 = (A, B, C) \cdot \mathbf{e}_{qi}$, and · represents a dot product. The coordinates (pos_x, pos_y, pos_z) represent the 3D position of the voxel without translation. The boundary surface inside the voxel is defined by Ax + By + Cz + D = 0, where (A, B, C) is a normalized plane normal pointing from the solid region to the liquid region. The value $\mathbf{e}_{qi}$ is the velocity vector associated with the distribution in the red channel. The third pass is similar to the second pass, except that this time the blue and alpha channels are modified, and $\mathbf{e}_{qi}$ is the velocity vector corresponding to the blue channel distribution.

In this way, all the boundary nodes are covered by the voxels. We then compute D at the beginning of the boundary condition fragment program with the following equations:

$$\Delta' = flag_1 / flag_2,$$

$$\Delta = 1 - \Delta'.$$

The meanings of $flag_1$ and $flag_2$ are the same as before. Note that each channel computes its $flag_1$, $flag_2$, and D independently, with its own $\mathbf{e}_{qi}$. A distribution is a boundary distribution only if, for the corresponding color channel, 1 ≥ D ≤ 0. If it is not a boundary distribution, the fragment program prevents modifying the corresponding color channel by assigning it the old value.

47.5 Visualization

To visualize the simulation, we inject particles into the flow field. The positions of the particles are stored in a texture and are updated every time step by a fragment program using the current velocity field. Similar to the generation of boundary voxels, the updated particle positions are then copied to a vertex array residing in the graphics memory for rendering. The whole simulation and rendering cycle is inside the GPU, hence there is no need to transfer large chunks of data between the main memory and the graphics memory. To better display the flow field, we arrange particles in a regular grid before injection and color them according to the position where they enter the flow field, as shown in Figure 47-1. Due to the requirements of the vertex array, the total number of particles is constant during the simulation. We use a fragment program to recycle particles that flow out of the border of the field or stay in a zero-velocity region and place them back at the inlet. If two particles coincide at exactly the same location at the same time, they will never separate during the advection. Visually, we will see fewer and fewer particles. To avoid this, we add a random offset to each particle when placing it at the inlet.

In 2D LBM, there is only one velocity slice, while in 3D LBM, the velocities form a volume. The advection fragment program fetches the velocity indicated by the current position of the particles. Therefore, the fragment program needs to access either a 3D texture or a 2D texture storing the flat volume. We chose the flat volume storage, which is much faster. Please note that if we stored the velocity as a stack of separate 2D textures, the advection would be very difficult.

47.6 Experimental Results

We have experimented with our GPU-based LBM implemented using OpenGL and Cg on an NVIDIA GeForce FX 5900 Ultra. This GPU has 256 MB of 425 MHz DDR SDRAM, and its core speed is 400 MHz. The host computer has a Pentium 4 2.53 GHz CPU with 1 GB PC800 RDRAM. All the results related to the GPU are based on 32-bit single-precision computation of the fragment pipeline. For comparison, we have also implemented a software version of the LBM simulation using single-precision floating point, and we measured its performance on the same machine.

Figure 47-8 shows a 2D flow field simulation based on the D2Q9 model with a lattice size of 256^2. We insert two circles (shown in red) into the field as the obstacles' boundary conditions. Vortices are generated because of the obstacles. The figure shows only a snapshot of the flow; the flow is relatively stable, but not constant. Figure 47-1 shows 3D flow simulations using the D3Q19 model with a lattice size of 50^3. Note that our simulation can handle arbitrarily complex and dynamic boundaries, which usually results in a complex flow field. We show only particles injected from a slit, to reduce clutter. The motion of the particles and the ability to change the viewing angle help in understanding the shape of the flow.

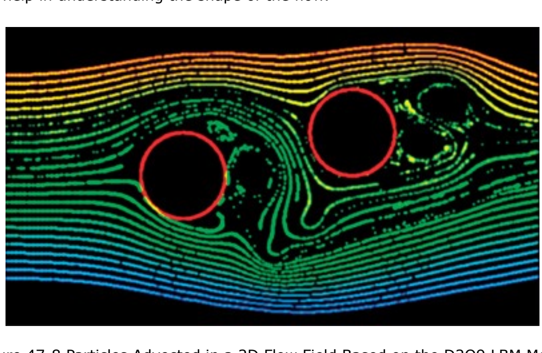

Figure 47-8 Particles Advected in a 2D Flow Field Based on the D2Q9 LBM Model

Figure 47-9 shows the time in milliseconds per step of the 3D LBM model as a function of the lattice size, running on both the CPU and the GPU. Note that both the *x* and *y* axes are in logarithmic scale. Figure 47-10 compares the two from a different perspective by showing the speedup factor. The time includes both simulation and visualization. Note that there is no need to transfer the velocity field or the particle positions between the main memory and the graphics memory. The time spent on advecting and rendering the particles is negligible with respect to the simulation. The speedup factor varies between 8 and 9 for most of the lattice sizes. When the lattice size approaches 128^3, the speedup is as high as 15. We are sure to get higher speedups with more recent GPUs, such as a GeForce 6800-class GPU. The step in the GPU timing curve—as well as in the GPU versus CPU speedup—is good evidence showing their different cache behavior. When the grid size gets bigger and surpasses a certain threshold, cache misses significantly slow down the CPU version. Due to the mostly sequential access patterns of the LBM algorithm and the GPU's optimization for sequential access, the cache-miss rate on the GPU is relatively independent of the grid size, and we don't see such a performance jump. The performance of the CPU version actually drops dramatically for lattice size of exactly power of two, possibly caused by 64K aliasing (Intel 2004). Therefore, we actually use lattices of size power-of-two-plus-one for the measurement. The CPU-based LBM implementation does not utilize SSE.

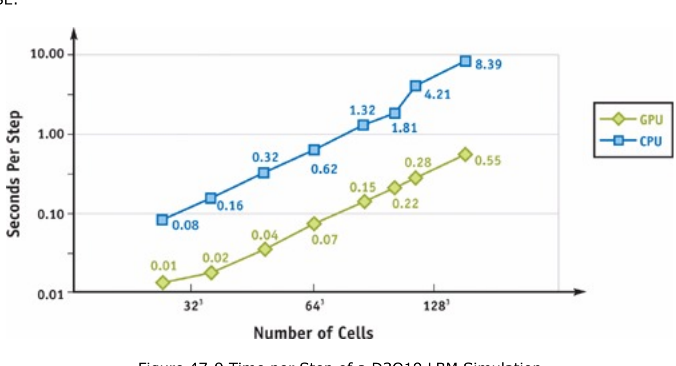

Figure 47-9 Time per Step of a D3Q19 LBM Simulation

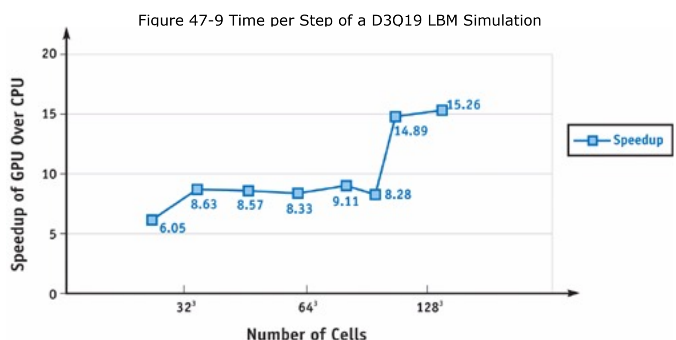

Figure 47-10 Speedup of the LBM on the GPU

47.7 Conclusion

We have presented a fluid flow simulation using the physically based Lattice Boltzmann Method, accelerated by programmable graphics hardware. The LBM simulation and its boundary conditions are of second-order accuracy in both time and space. Our GPU-based simulation using floating-point computation achieves the same accuracy as the CPU-based LBM simulation, yet it is much faster. Our experimental results have shown that the GPU version is significantly faster (for example, 8 to 15 times faster, for the D3Q19 model) for most lattice sizes in both 2D and 3D simulations. To attain these speeds, we have incorporated several optimization techniques into the GPU method. We have also proposed a GPU-based general-purpose voxelization algorithm and its extension to handle arbitrarily complex and dynamic boundary conditions in real time.

47.8 References

Other graphics researchers have explored the approach of solving the Navier-Stokes (NS) equations to simulate amorphous phenomena, such as gases and fluid: Foster and Metaxas 1997, Stam 1999, and Fedkiw et al. 2001. Readers may find a comparison of direct NS solvers with the LBM in Wei et al. 2004. Researchers have also developed GPU-accelerated conjugate gradient solvers that can be used for NS equations: Bolz et al. 2003, Krüger and Westermann 2003. There are also other works on using GPUs to compute fluid dynamics: Harris et al. 2003, Goodnight et al. 2003, and Harris 2004. More details of the GPU-accelerated LBM can be found in Li 2004.

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Dedication

To everyone striving to make today's best computer graphics look primitive tomorrow