Implementation of Reynolds Equation Solver on GPGPU for Gas Film Lubrication Problem Ji-Hoon Kang, Hun Joo Myung, Kwang Jin Oh, Chan Yeol Park, Supercomputing Center, KISTI



Implementation on GPGPU

Gauss-Siedel iteration

Reynolds eqn. for gas film pressure

 $a_{P}P_{P} = a_{E}P_{E} + a_{W}P_{W} + a_{N}P_{N} + a_{S}P_{S} + bP_{P}^{(n-1)}$

Mass continuity eqn. for pocket pressure

 $\sum_{f_{Pocket}} a_P P_{Pocket} = \sum_{f_{Pocket}} \left(a_E P_E + a_W P_W + a_N P_N + a_S P_S + b P_{Pocket}^{(n-1)} \right)$ $+A_{Supply}P_{Supply}$

Red/Black Gauss-Siedel iteration (RBGS)

In iteration n:

► <u>1st pass</u> All red nodes are updated using old values of black nodes.

2nd pass

All black nodes are updated using updated values of red nodes

No data dependency among nodes of same colors

\rightarrow the nodes can be assigned to each thread and updated in parallel within each pass

Assign each z node to each thread and each theta node to each block

// Red/Black Gauss-Siedel iterative pressure solver structure void Pressure_CUDA::iterate(){ int redblack: unsigned int threads = NzNode; unsigned int blocks = NthNode; redblack = 0; //1st pass iteratePressure_kernel<<<blocks, threads, Mem_size>>>(d_MeshType,...,redblack); cutilSafeCall(cudaMemcpy(err,d_err, sizeof(double)*blocks, cudaMemcpyDeviceToHost)); for(i=0;i<blocks;i++) { error+=err[i]; } // Error summation</pre> redblack = 1; **//2nd pass** iteratePressure_kernel<<<blocks, threads, Mem_size>>>(d_MeshType,...,redblack); cutilSafeCall(cudaMemcpy(err,d_err, sizeof(double)*blocks, cudaMemcpyDeviceToHost)); for(i=0;i<blocks;i++) { error+=err[i]; } // Error summation</pre> }while(error>criteria); // Computational kernel structure _global__ void iteratePressure_kernel(int *d_MeshType, …, int redblack) { int ildx = threadldx.x: int jldx = blockldx.x; if(((ildx+jldx)&1) == redblack && d_MeshType[jldx*const_NzNode+ildx]==-1) { ... Coefficient calculation aP=aE+aW+aN+aS+Fe-Fw+b; d_dP[jldx*const_NzNode+ildx]=(aE*PE+aW*PW+aN*PN+aS*PS+b*oldPP)/aP-PP; **d_P[jldx*const_NzNode+ildx]** +=0.4*d_dP[jldx*const_NzNode+ ildx]; // ... Error calculation

double_reduction(error, d_err);







Wall clock time comparison



|112x160|-CPU+GPU 16.9 CPU 3074.5 252x360 CPU+GPU 256.2 26.8 (10%) 33.4 (13%) 54.1(21%) 141.9 (55%)

A small serial part becomes a critical bottleneck as the problem size gets bigger and GPU speedup increases. Not only core calculation part but also general calculation part needs to be parallelized for better speedup.

Conclusion & Future work

- lubrication problem.
- size gets bigger and GPU efficiency increases.

Mesh size 112x160		Mesh size 252x360	
50		3000	
00		2 2000	
50 — —	8.7x		12.0x
0		S 0	
CPU only CPU+GPU CPU only CPU+GPU			
Others 🛛	MemCopy	Others	MemCopy
teratePocketP	IterateP	IteratePocketF	P 🗖 IterateP
	_		
IterateP(s)	ItrPocketP(s)	MemCp(s)	Others(s)
17.7 (86%)	2.2 (11%)	-	0.7 (3%)
0.8 (16%)	2.6 (52%)	0.8(15%)	0.8 (16%)
135.4 (92%)	6.0 (4%)	-	5.7 (4%)
2.3 (14%)	4.9 (29%)	3.7(22%)	6.0 (36%)
2961 6(02%)			

Implemented a Reynolds equation solver on GPGPU for gas film

 Achieved 106x speedup for core calculation part and overall 12x speedup (DP) using RBGS, relative to 1 core of AMD Llano A8-3850. • A tiny serial part can become a critical bottleneck as the problem

• Future work will include the development of general gas film analysis solver and the development of parallelization scheme for remaining serial part, such as integration, error check, and et al.