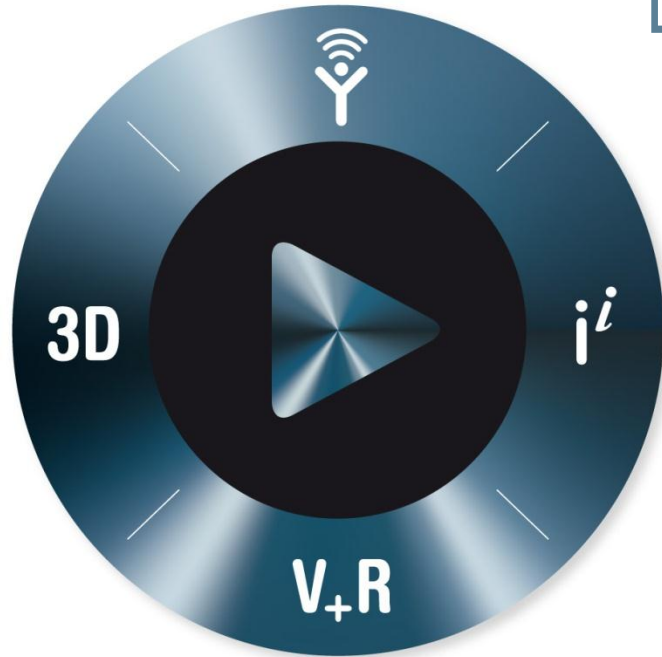


# Evolving Use of GPU for Dassault Systemes Simulation Products



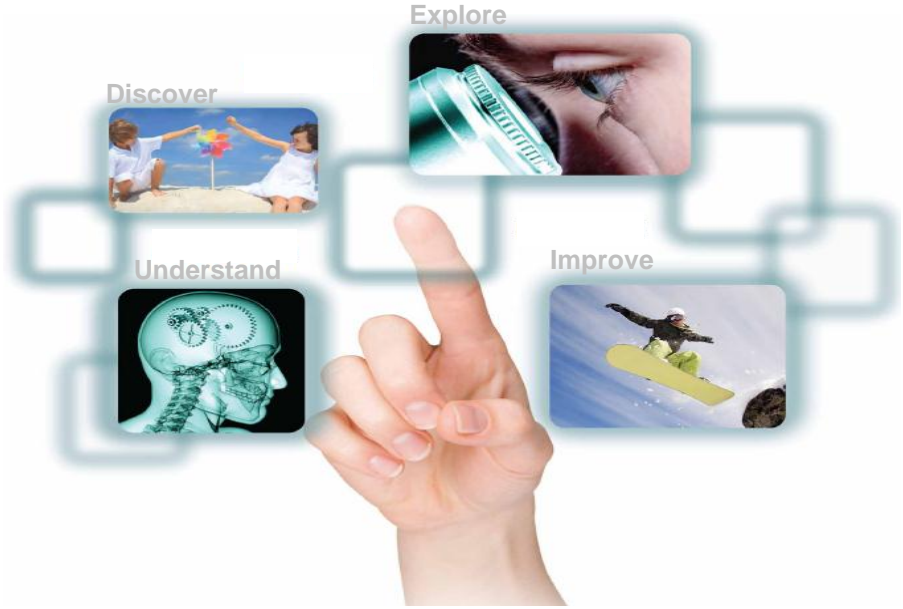
**3D**EXPERIENCE

Luis Crivelli  
And  
Matt Dunbar

# Dassault Systèmes is dedicated to making...

Realistic Simulation  
*an integral*  
business practice

to Explore,  
Discover,  
Understand,  
Improve

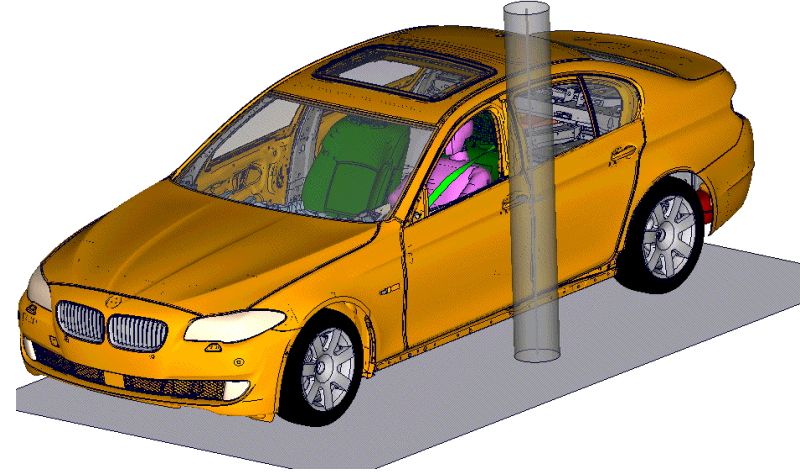


## *product, life, & nature*

# Simulation with Abaqus

## Finite Element Simulation

Abaqus/Standard – static structural simulations  
Abaqus/Explicit – short term dynamic simulations



"Predictive Crashworthiness Simulation in a Virtual Design Process without Hardware Testing",  
Jurgen Lescheticky, Hariakto Hooputra and Doris Ruckdeschel, BMW Group, SIMULIA  
Customer Conference, May 2010

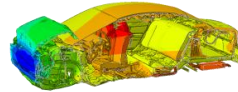
# Simulation runtimes

- ▶ Simulation is a valuable part of an engineering design process, but computational cost is significant



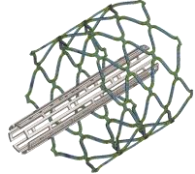
## Bottle Stacking

- Routine static analysis
- Compute time of 4-6 hours on 4 x86 cores



## Automotive Crash

- Complex dynamics model
- Compute time of 2-5 days using between 16 and 32 x86 cores



## Stent Expansion

- Complex static analysis
- Compute time of 12-24 hours using between 8 and 32 x86 cores

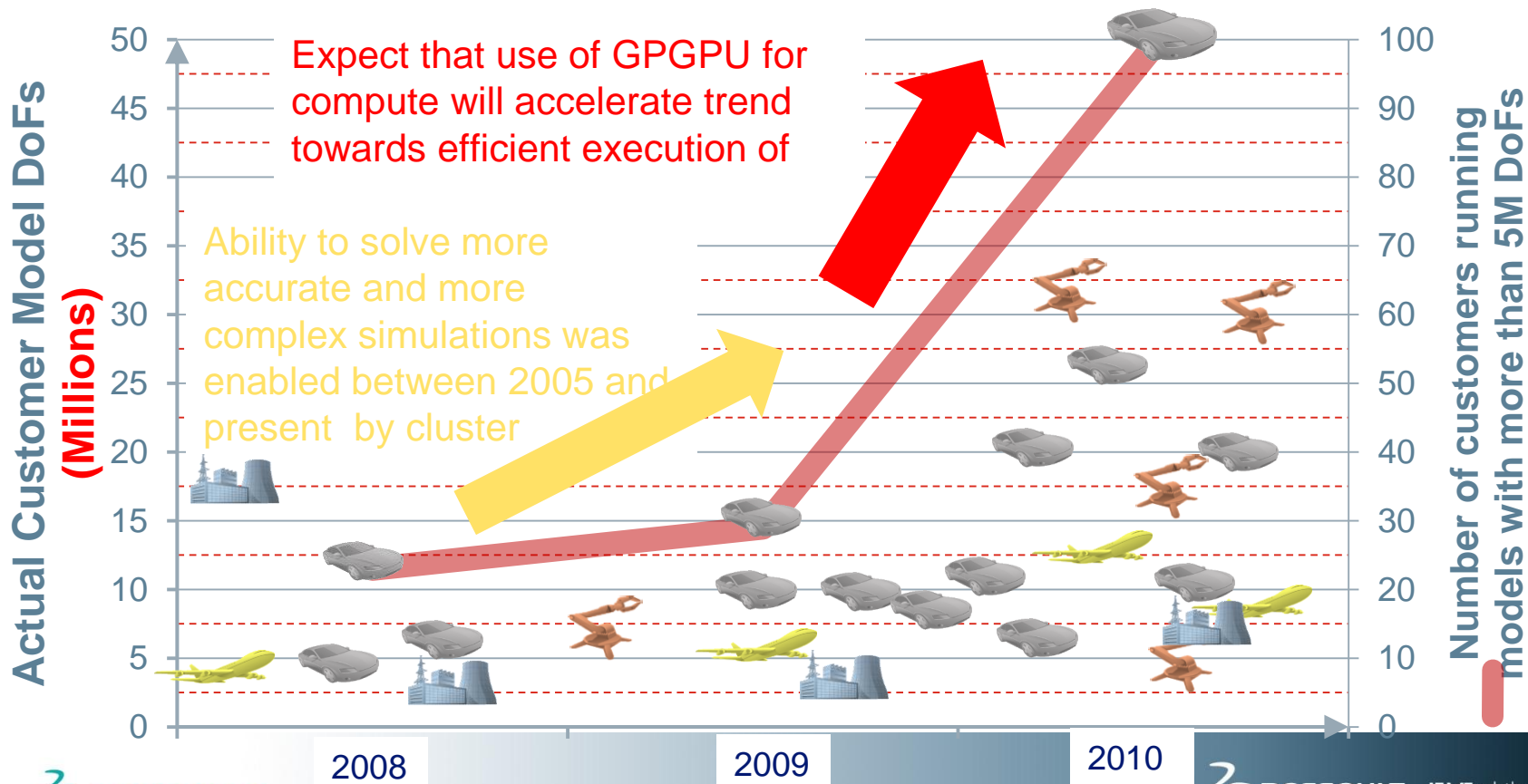


## Gasket Sealing

- Large static model
- Compute time of 2-5 days using between 32 and 64 x86 cores

**Reducing compute cost is critical!**

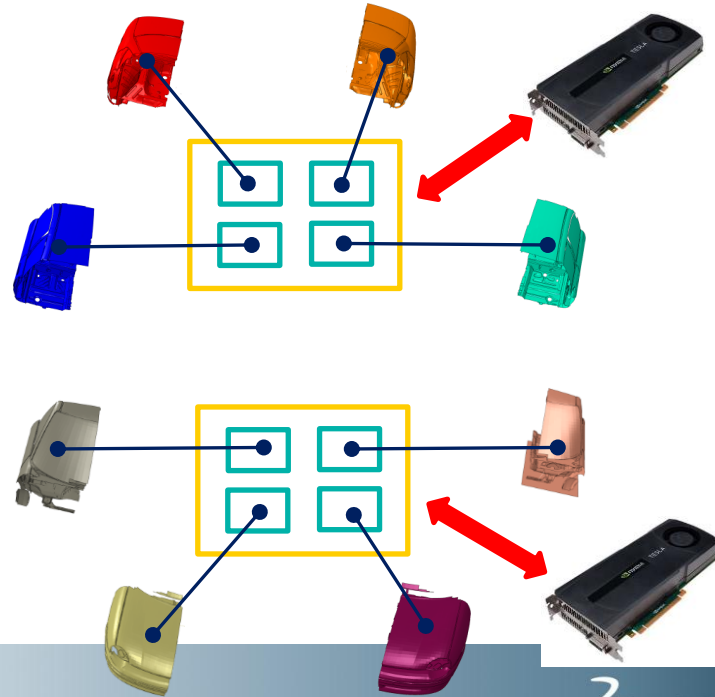
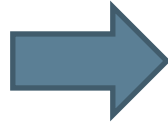
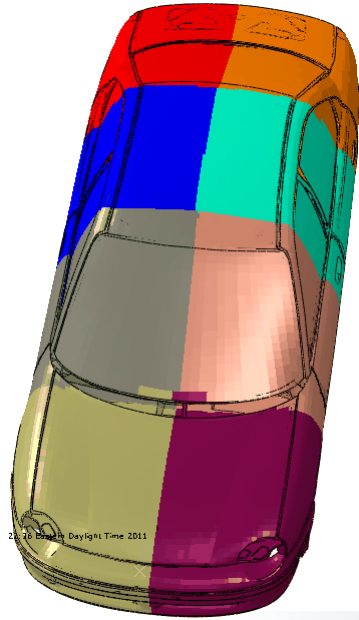
# Demand for Higher Accuracy



3DS.COM/SIMULIA © Dassault Systèmes | Confidential Information | 5/16/2012 | ref.: 3DS\_Document\_2012

# Simulation compute architecture

- ▶ Computational work is distributed to many cores on multiple servers by splitting a model into domains which are assigned to cores to parallelize computations

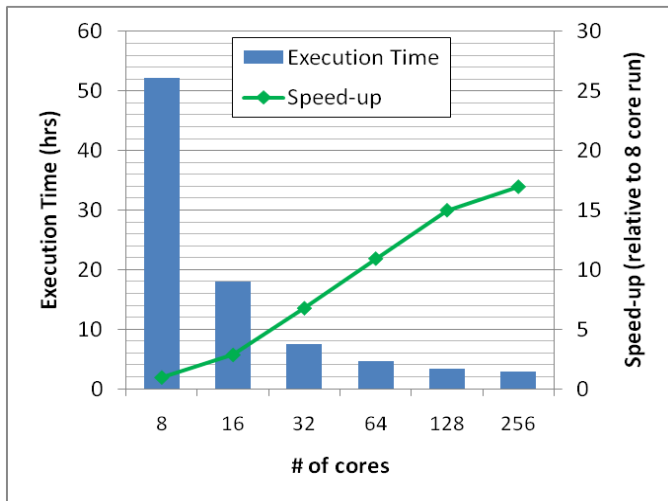


**New code allows x86 cores to offload key compute work onto one or more Nvidia Tesla cards to accelerate computation**

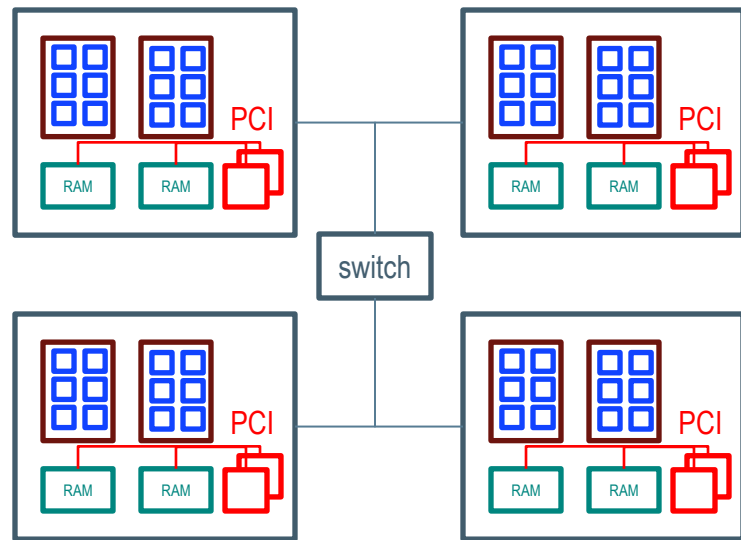
# Abaqus

## Existing Cluster-based Architecture

Abaqus applications allow users to exploit X86 clusters to decrease runtimes

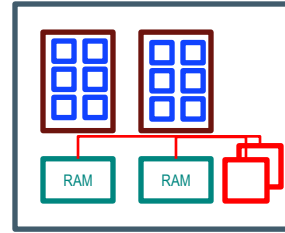


Can GPU deliver a faster or more efficient solution?



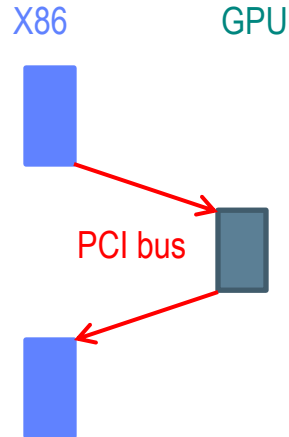
Machine  
Socket  
Core  
GPU

# Approaches to Exploiting GPU



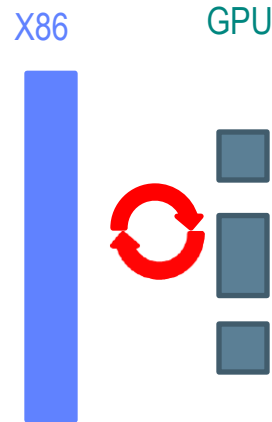
## Offload Code to Card

- X86 cores are idle while GPU card processes



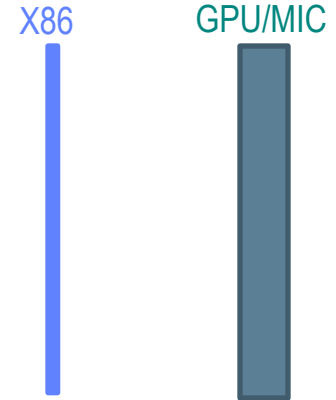
## Hybrid Mode

- X86 cores and GPU card are used simultaneously with X86 assigning appropriate work to GPU



## GPU As Platform

- Very limited "control" compute on X86; bulk of work done on GPU card





# Abaqus Solvers

## Key Code Components

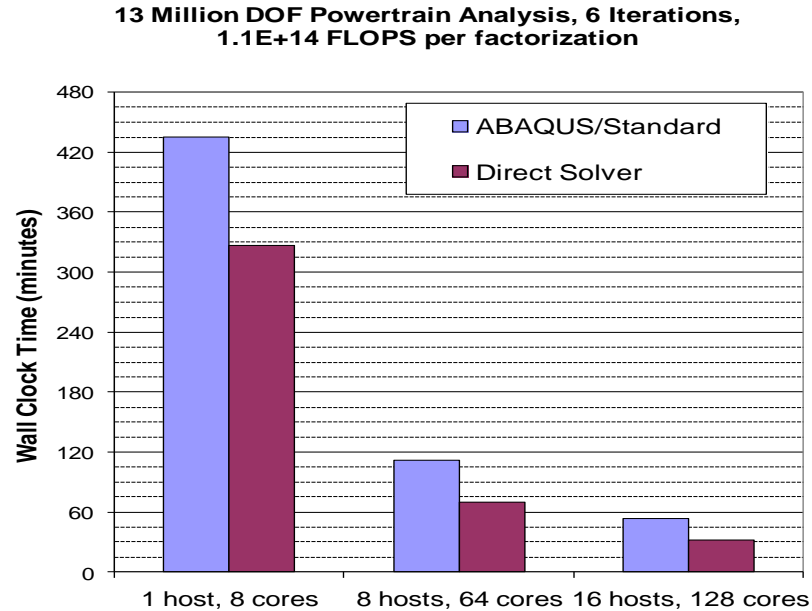
	Code Component	Cost of Code	Nature of Code
Abaqus/Standard	Linear Equation Solver	Increases with problem size.	Dense linear algebra kernels and “control code.” Relatively small amount of code with high computational cost.
	Finite Elements	Most significant cost other than equation solver.	Naturally parallel but code is not written to expose SIMD parallelism.
Abaqus/Explicit	Elements	Typically 50%-75% of cost.	Naturally parallel. Code is written for SIMD architectures.
	Constraints and other	Much of remaining cost.	Complex parallelism.

Uses GPU in Abaqus 6.11 and 6.12

Focus of prototyping work

# Abaqus/Standard

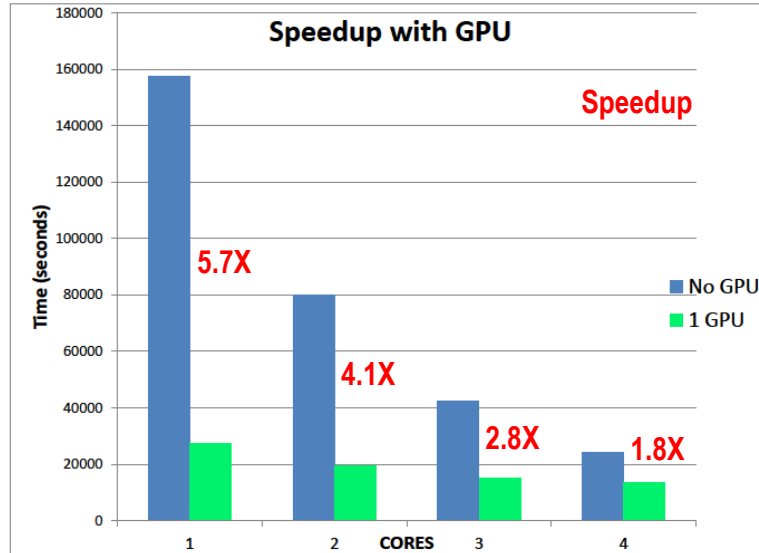
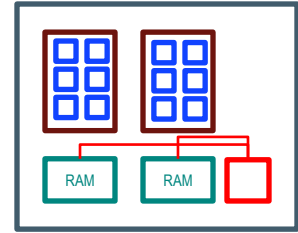
## Time in Equation Solver



- ▶ As noted on previous slide Abaqus/Standard equation solver is a significant cost in execution of a simulation
- ▶ Natural target for GPU acceleration

# Abaqus/Standard

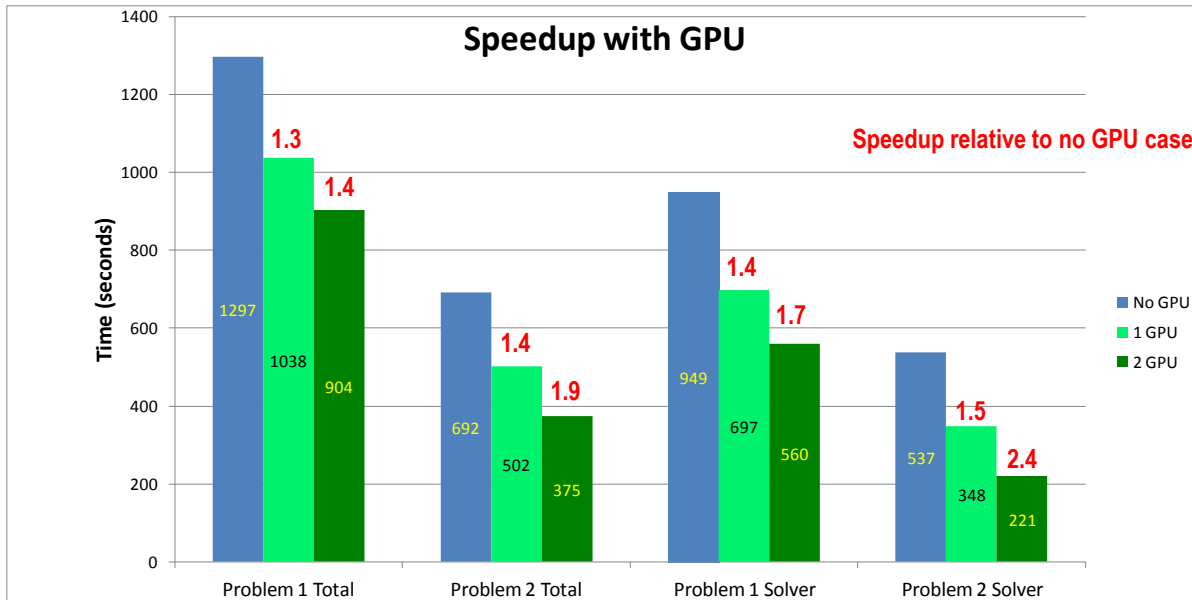
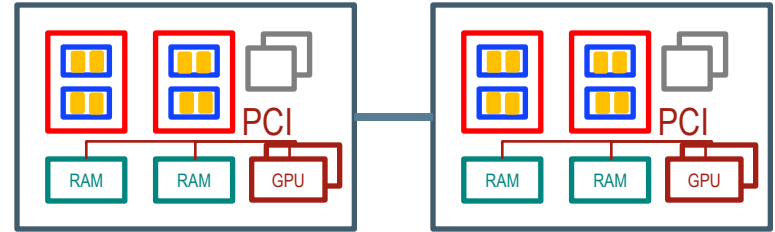
## Single GPU/Single Server



- ▶ At lower core counts GPU is more effective because the accelerated code takes a higher percentage of the overall time (code accelerated by the GPU is also effectively parallelized on X86 cores)
- ▶ Beyond 4 cores, X86 processors start overwhelming the GPU with computation

# Abaqus/Standard

Cluster and multi-GPU support – hybrid mode



Jobs run on 2 hosts each with 12 cores, 48 GB of RAM, and 2 Nvidia GPU's

Problem 1

- 1.5 MDOF
- 5.37 teraflops per iteration

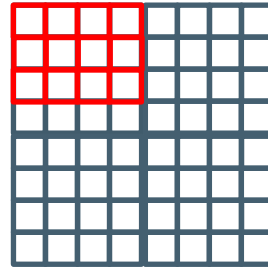
Problem 2

- 3.4 MDOF
- 10.8 teraflops per iteration

# Abaqus/Explicit

## GPU as a platform using Portland Group (OpenACC) compiler

- ▶ Explicit finite element codes do not have a natural bottleneck
  - ▷ Compute cost is spread through many 100's of routines
  - ▷ Code does have a natural SIMD structure
- ▶ SIMULIA has done a prototype in which the Portland Group compiler was used to build existing X86 code for GPU



Process group 1

Loop over groups of elements

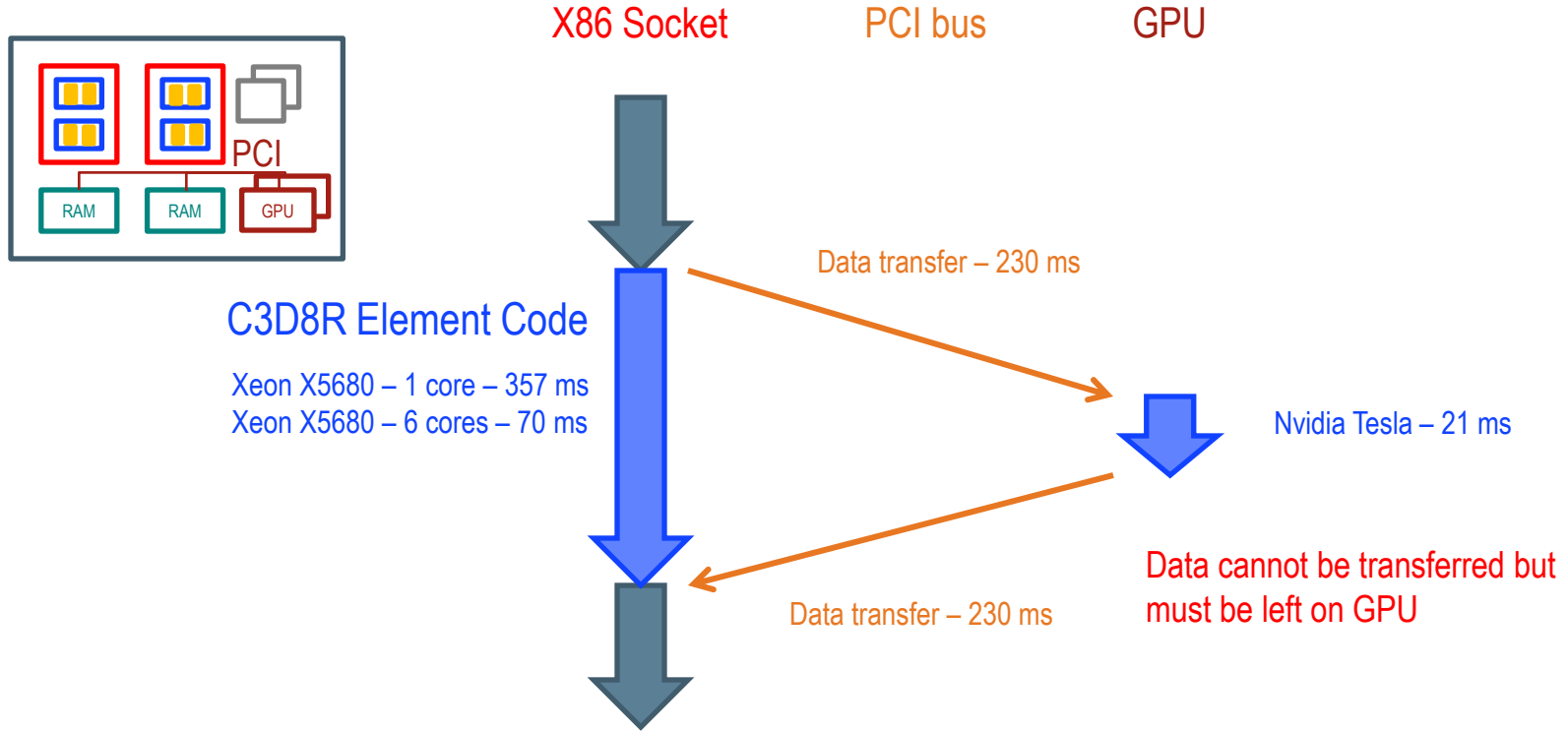
Loop A – process 1 to ngroup elements

Loop B – process 1 to ngroup elements

Loop C – process 1 to ngroup elements

Parallelization is done at a fine-grained level. Generally relies on elements being processed being identical (Single Instruction Multiple Data)

# Abaqus/Explicit Prototype



# Memory bandwidth bounds GPU performance

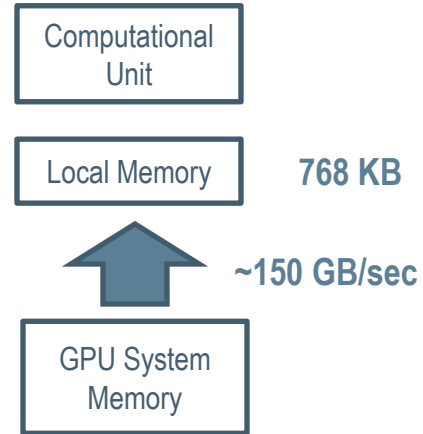
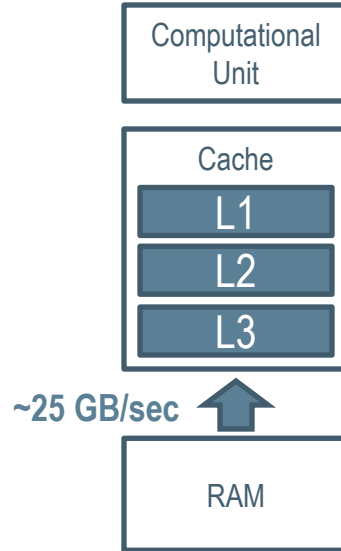
- ▶ Given peak performance of GPU near 1 TFlop vs peak performance for Westmere of ~70 Gflops why does Explicit code sped up by only 3X?

On X86 system temporary arrays remain in cache, but on GPU limited local memory is shared between a large number of threads so temporary data ends up be written back to system memory.

```
do k = 1, groupsize
  temp1(k) = in1(k) * in2(k)

do k = 1, groupsize
  temp2(k) = temp1(k) * in3(k)

do k = 1, groupsize
  out(k) = temp2(k) * x
```



High degree of data parallelism, but each piece of data is used a small number of times in operations

# Conclusions

- ▶ Use of GPU as an X86 accelerator in a hybrid mode is effective and continued area of development
- ▶ Investigations into GPU as a platform are ongoing
  - ▷ Experience with OpenACC compiler approach has been good
  - ▷ Highly data parallel code does have bandwidth issues on GPU which limits gains
  - ▷ Data management between X86 memory and GPU is a key topic



