Introduction to MPP version of LS-DYNA®

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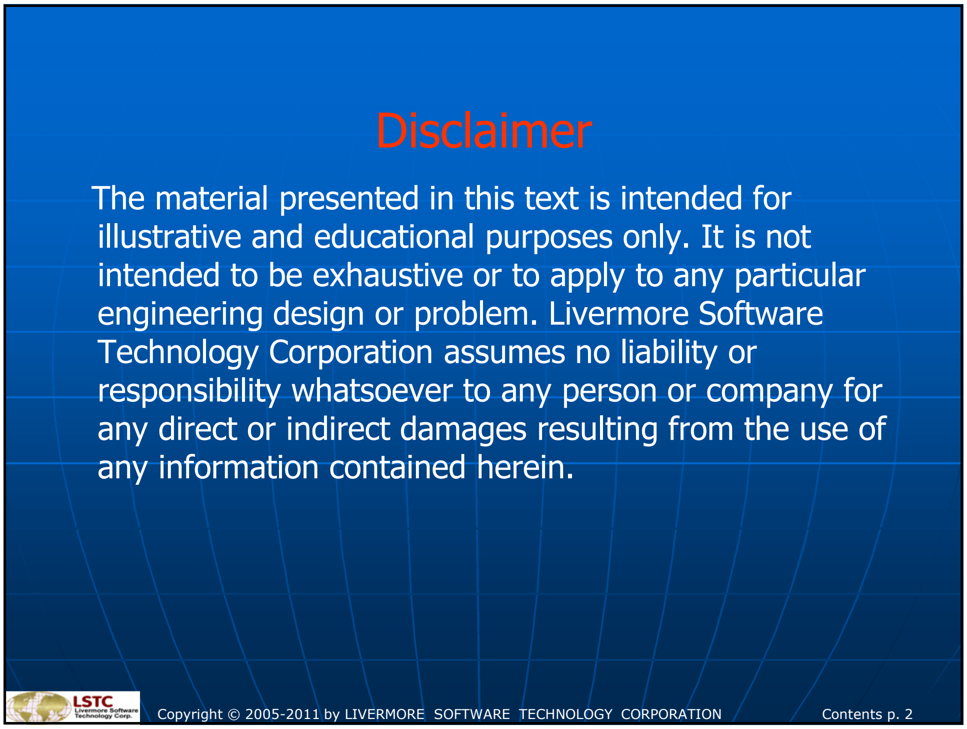
Livermore Software Technology Corporation



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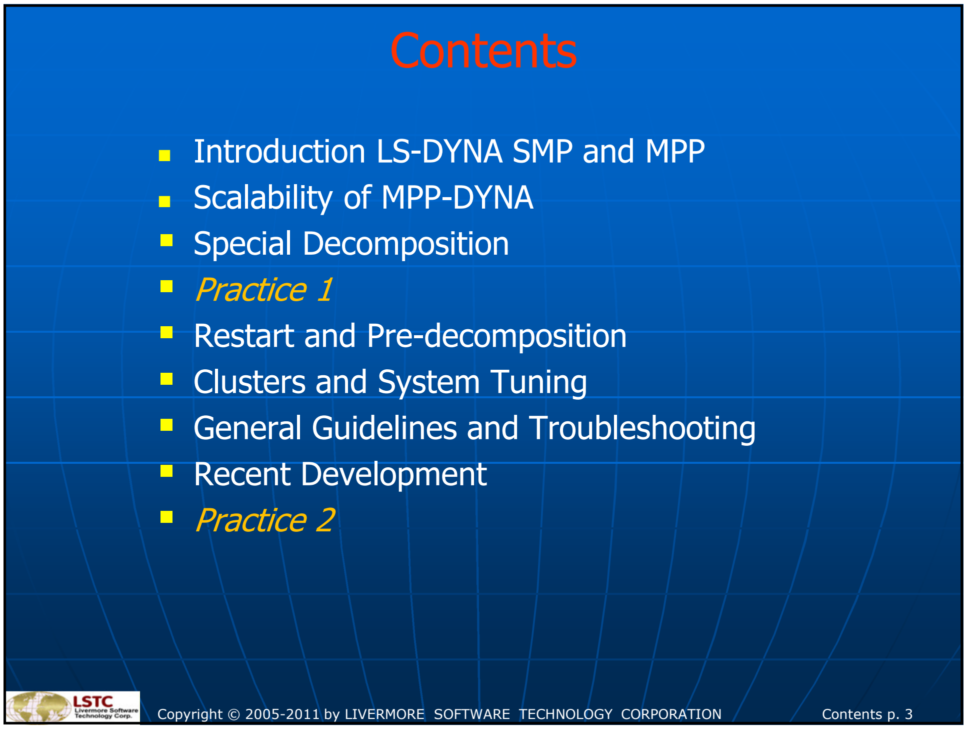

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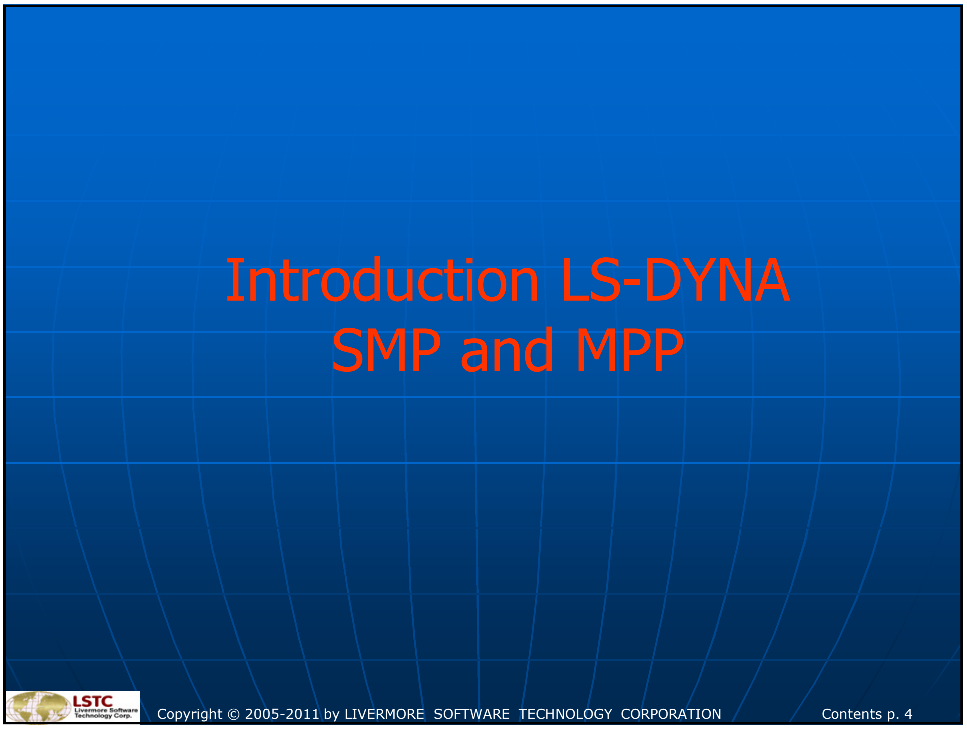
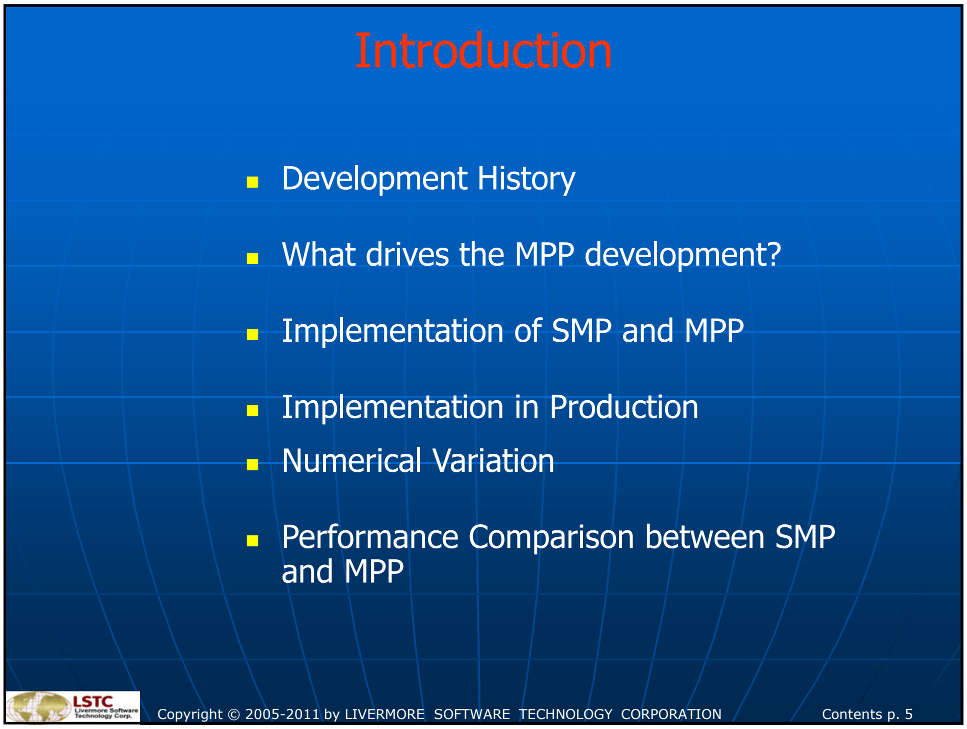
- Scalability of MPP-DYNA
- Special Decomposition
- *Practice 1*
- Restart and Pre-decomposition
- Clusters and System Tuning
- General Guidelines and Troubleshooting
- Recent Development
- *Practice 2*

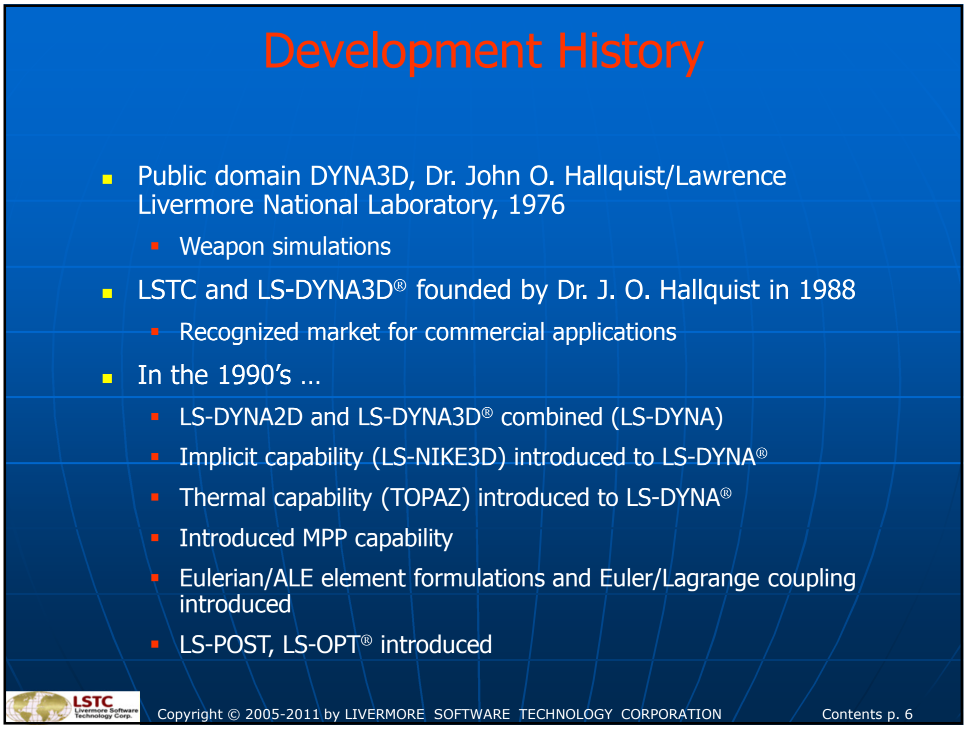


Introduction LS-DYNA SMP and MPP



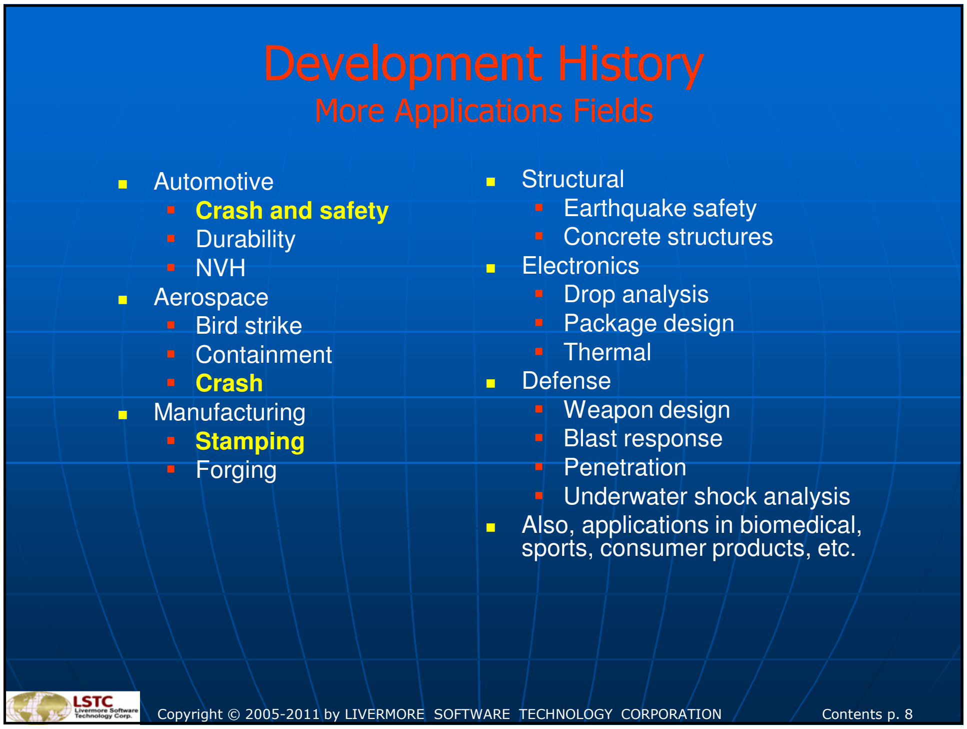
Introduction

- Development History
- What drives the MPP development?
- Implementation of SMP and MPP
- Implementation in Production
- Numerical Variation
- Performance Comparison between SMP and MPP



Development History

- Public domain DYNA3D, Dr. John O. Hallquist/Lawrence Livermore National Laboratory, 1976
  - Weapon simulations
- LSTC and LS-DYNA3D® founded by Dr. J. O. Hallquist in 1988
  - Recognized market for commercial applications
- In the 1990's …
  - LS-DYNA2D and LS-DYNA3D® combined (LS-DYNA)
  - Implicit capability (LS-NIKE3D) introduced to LS-DYNA®
  - Thermal capability (TOPAZ) introduced to LS-DYNA®
  - Introduced MPP capability
  - Eulerian/ALE element formulations and Euler/Lagrange coupling introduced
  - LS-POST, LS-OPT® introduced



Development History
More Applications Fields

- Automotive
  - **Crash and safety**
  - Durability
  - NVH
- Aerospace
  - Bird strike
  - Containment
  - **Crash**
- Manufacturing
  - **Stamping**
  - Forging
- Structural
  - Earthquake safety
  - Concrete structures
- Electronics
  - Drop analysis
  - Package design
  - Thermal
- Defense
  - Weapon design
  - Blast response
  - Penetration
  - Underwater shock analysis
- Also, applications in biomedical, sports, consumer products, etc.

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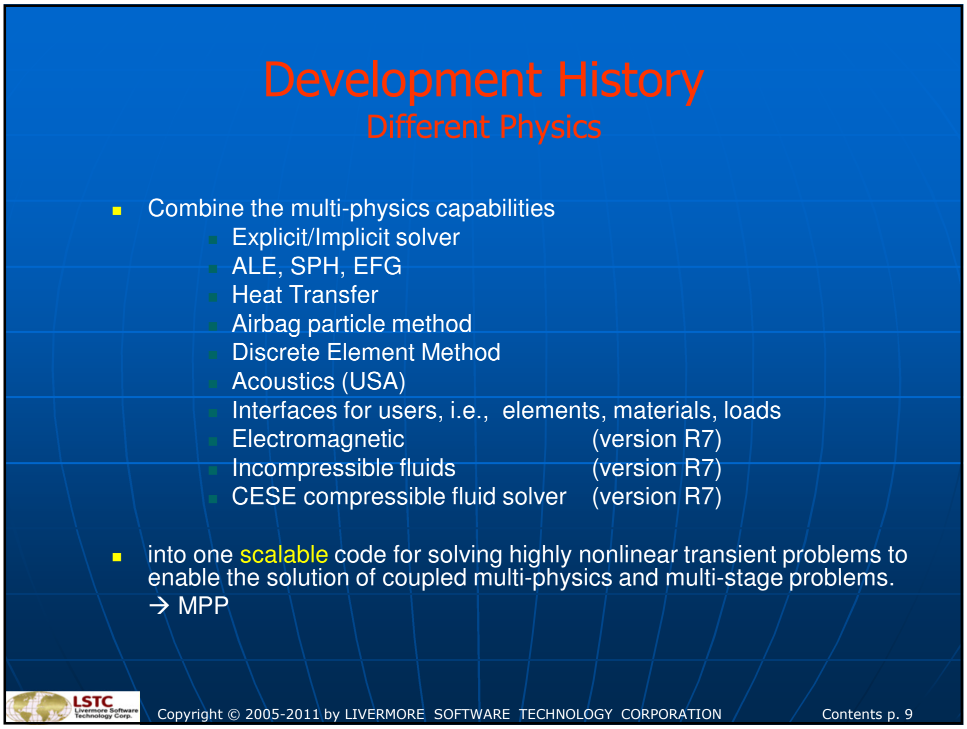

Development History
Different Physics

- Combine the multi-physics capabilities
  - Explicit/Implicit solver
  - ALE, SPH, EFG
  - Heat Transfer
  - Airbag particle method
  - Discrete Element Method
  - Acoustics (USA)
  - Interfaces for users, i.e., elements, materials, loads
  - Electromagnetic (version R7)
  - Incompressible fluids (version R7)
  - CESE compressible fluid solver (version R7)
- into one scalable code for solving highly nonlinear transient problems to enable the solution of coupled multi-physics and multi-stage problems. → MPP

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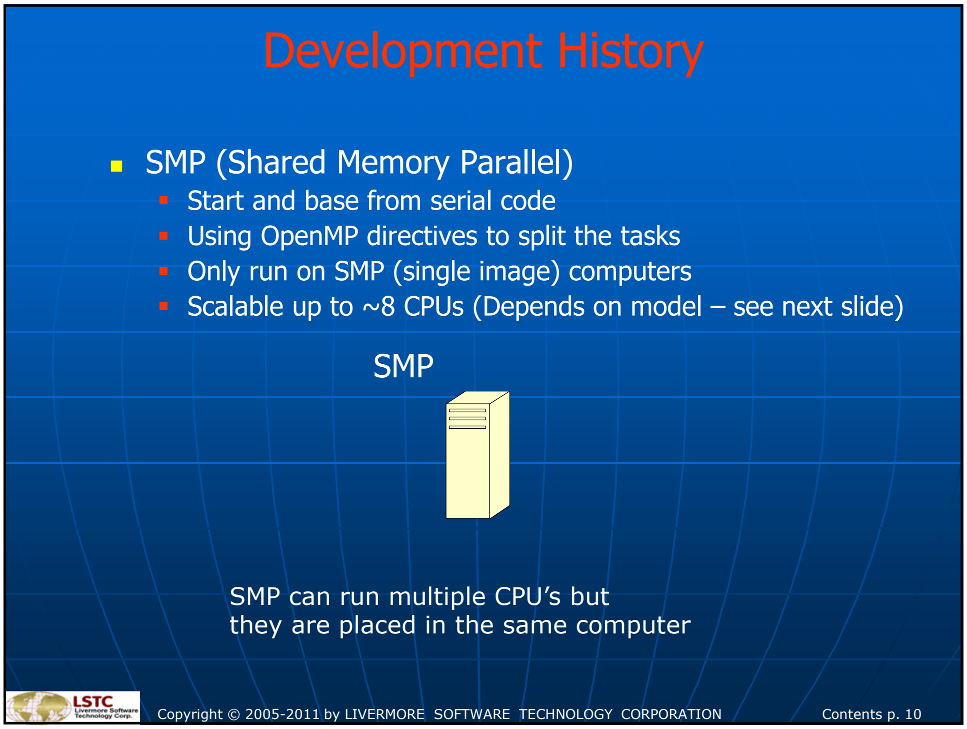


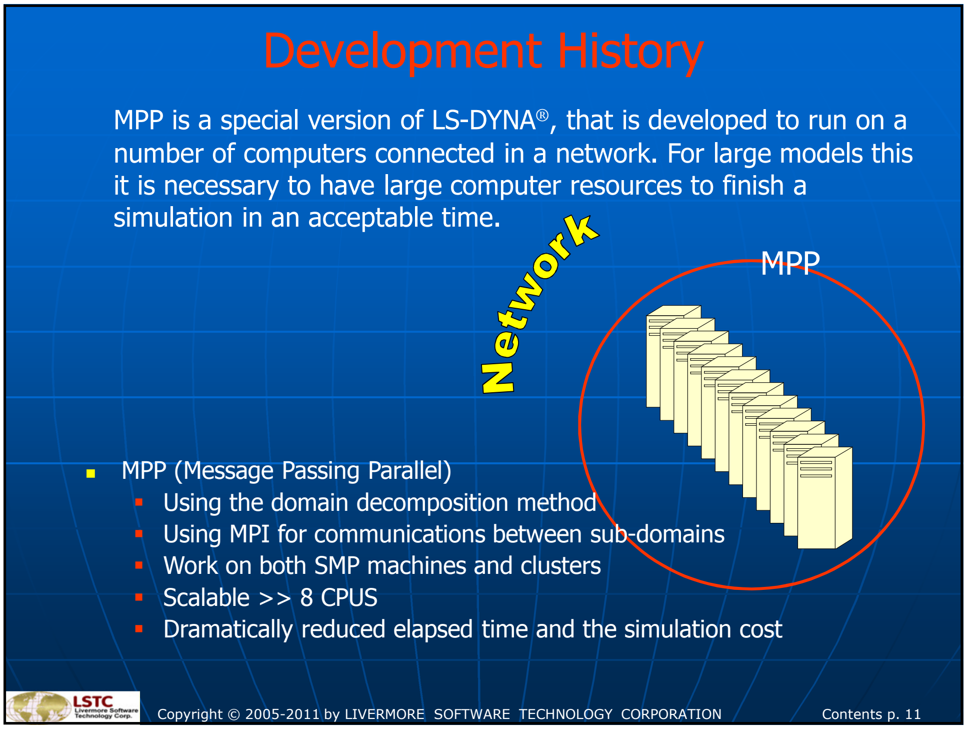
Development History

- SMP (Shared Memory Parallel)
  - Start and base from serial code
  - Using OpenMP directives to split the tasks
  - Only run on SMP (single image) computers
  - Scalable up to ~8 CPUs (Depends on model – see next slide)

SMP

SMP can run multiple CPU's but
they are placed in the same computer

LSTC
Livermore Software
Technology Corp.

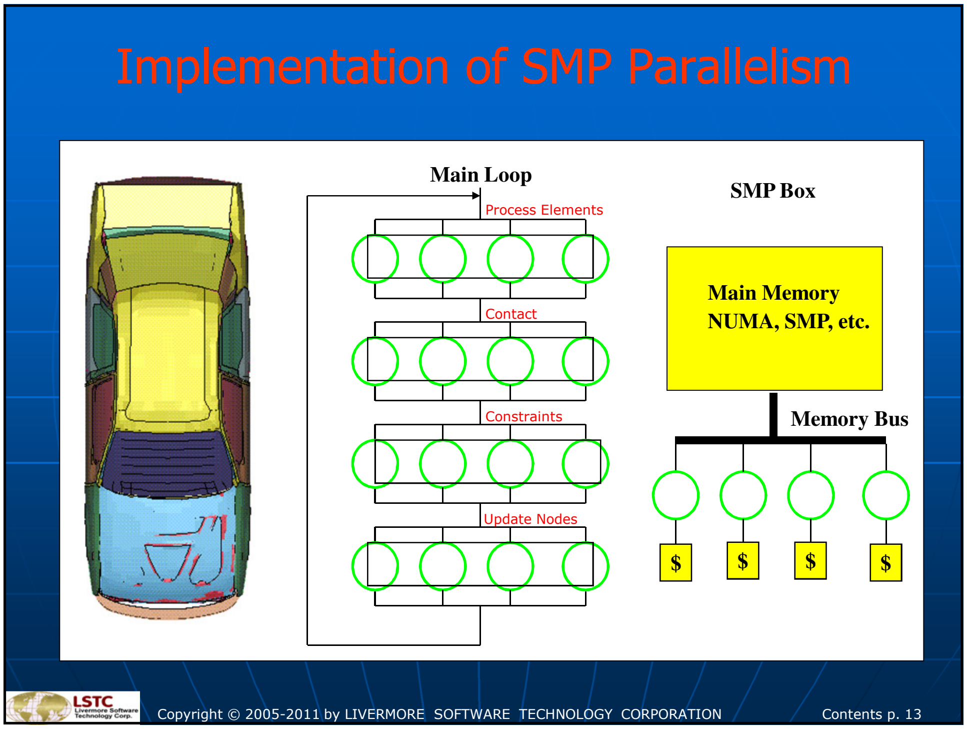


Development History

MPP is a special version of LS-DYNA®, that is developed to run on a number of computers connected in a network. For large models this it is necessary to have large computer resources to finish a simulation in an acceptable time.

Network
MPP

- MPP (Message Passing Parallel)
  - Using the domain decomposition method
  - Using MPI for communications between sub-domains
  - Work on both SMP machines and clusters
  - Scalable >> 8 CPUS
  - Dramatically reduced elapsed time and the simulation cost



Implementation of SMP Parallelism

Main Loop

Process Elements

Contact

Constraints

Update Nodes

SMP Box

Main Memory
NUMA, SMP, etc.

Memory Bus

$ $ $ $

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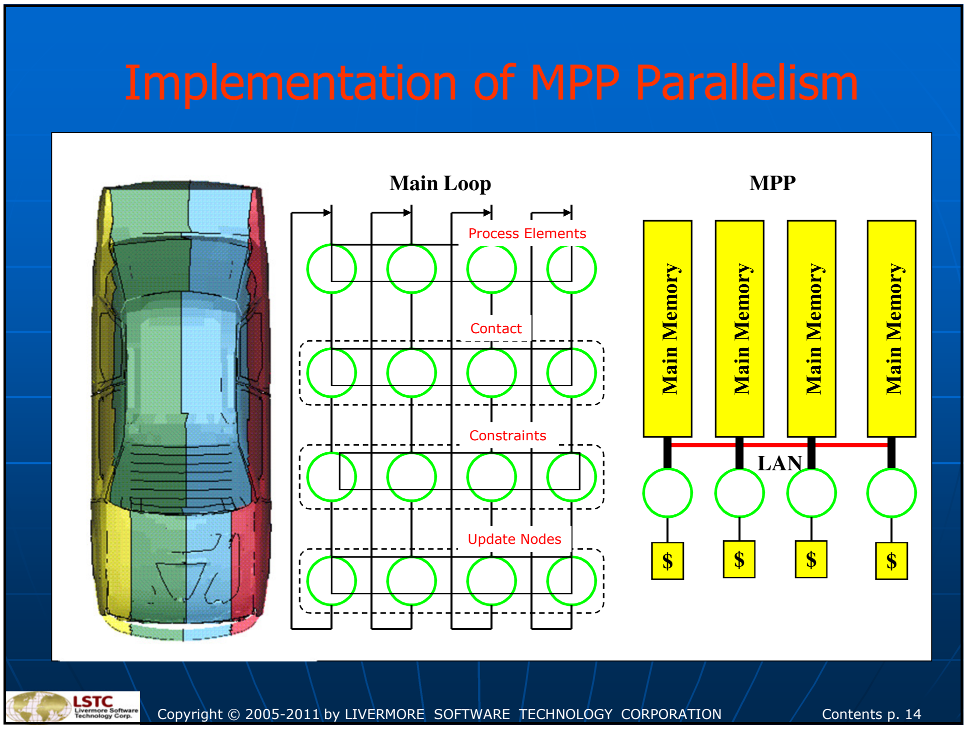

Implementation of MPP Parallelism

Main Loop

Process Elements

Contact

Constraints

Update Nodes

MPP

Main Memory

Main Memory

Main Memory

Main Memory

LAN

$ $ $ $

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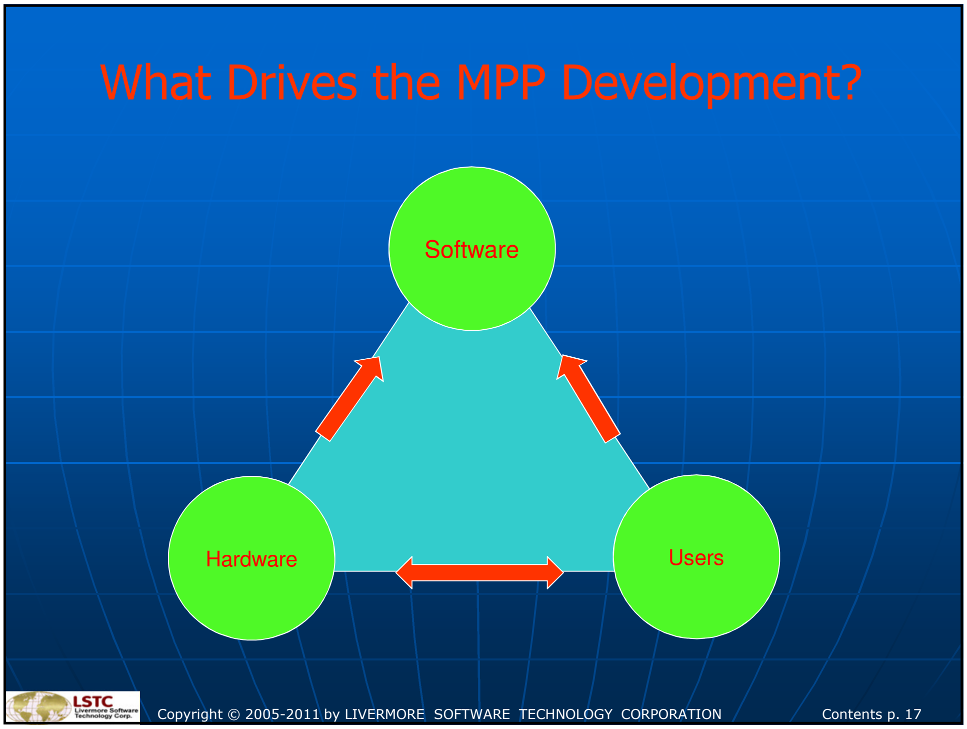
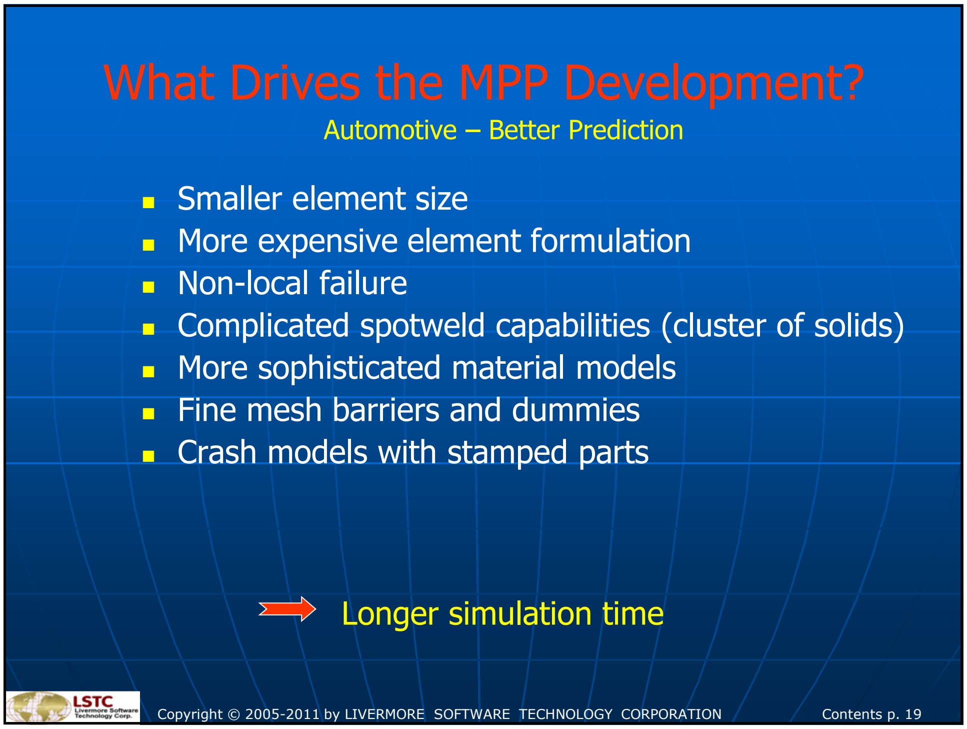


What Drives the MPP Development?

Software

Hardware

Users

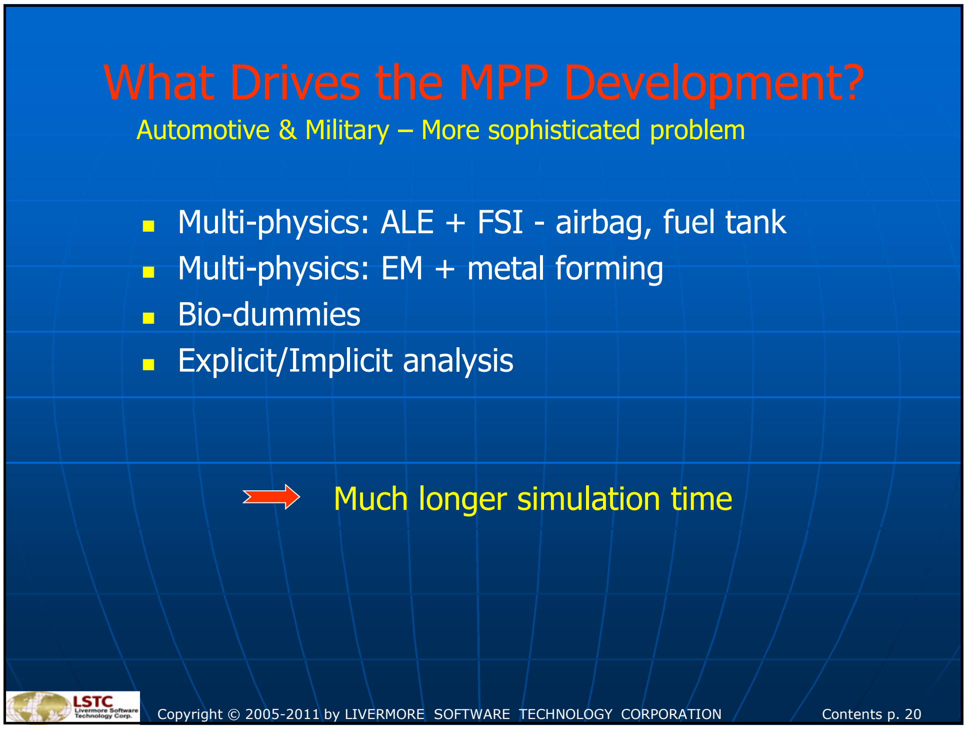


What Drives the MPP Development?

Automotive – Better Prediction

- Smaller element size
- More expensive element formulation
- Non-local failure
- Complicated spotweld capabilities (cluster of solids)
- More sophisticated material models
- Fine mesh barriers and dummies
- Crash models with stamped parts

→ Longer simulation time



What Drives the MPP Development?

Automotive & Military – More sophisticated problem

- Multi-physics: ALE + FSI - airbag, fuel tank
- Multi-physics: EM + metal forming
- Bio-dummies
- Explicit/Implicit analysis

→ Much longer simulation time

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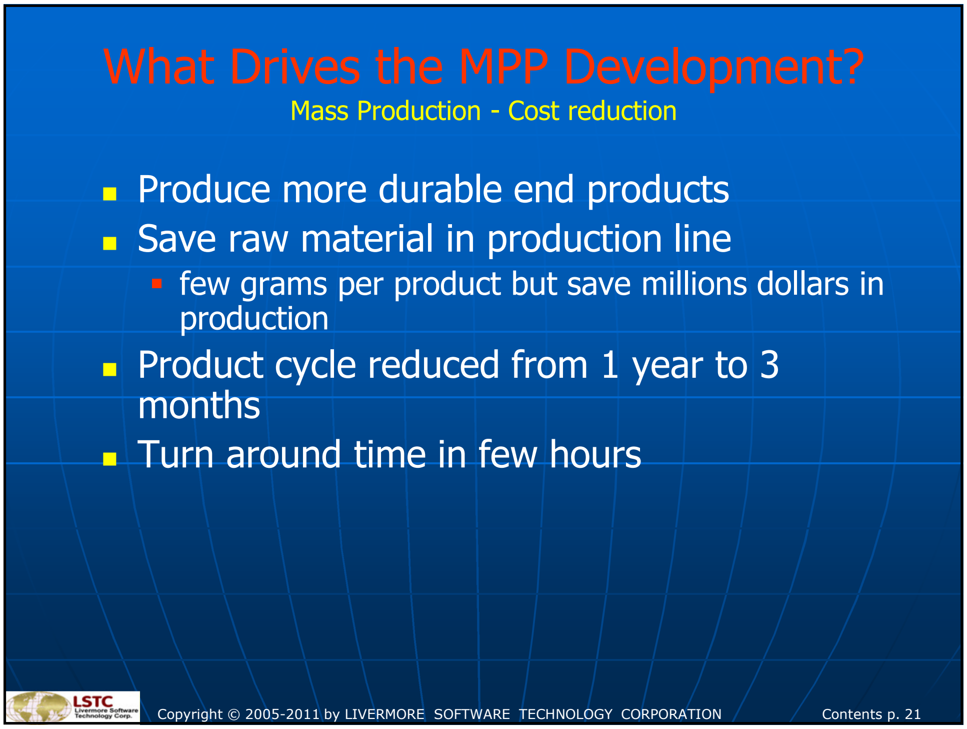

What Drives the MPP Development?

Mass Production - Cost reduction

- Produce more durable end products
- Save raw material in production line
  - few grams per product but save millions dollars in production
- Product cycle reduced from 1 year to 3 months
- Turn around time in few hours

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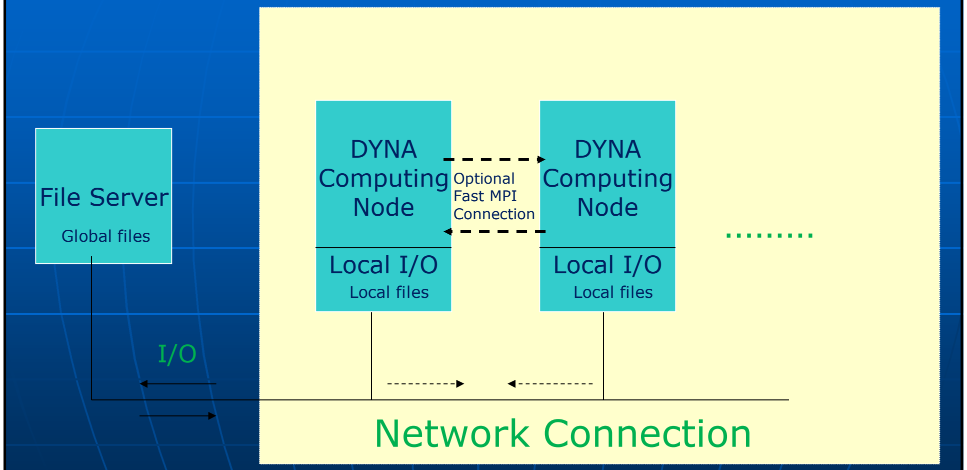
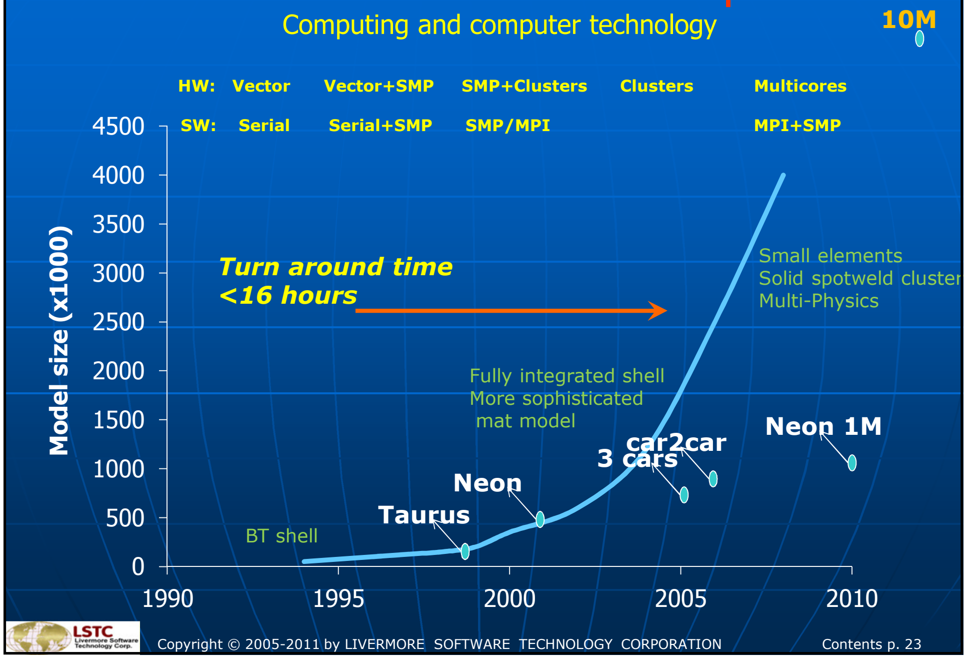
What Drives the MPP Development?

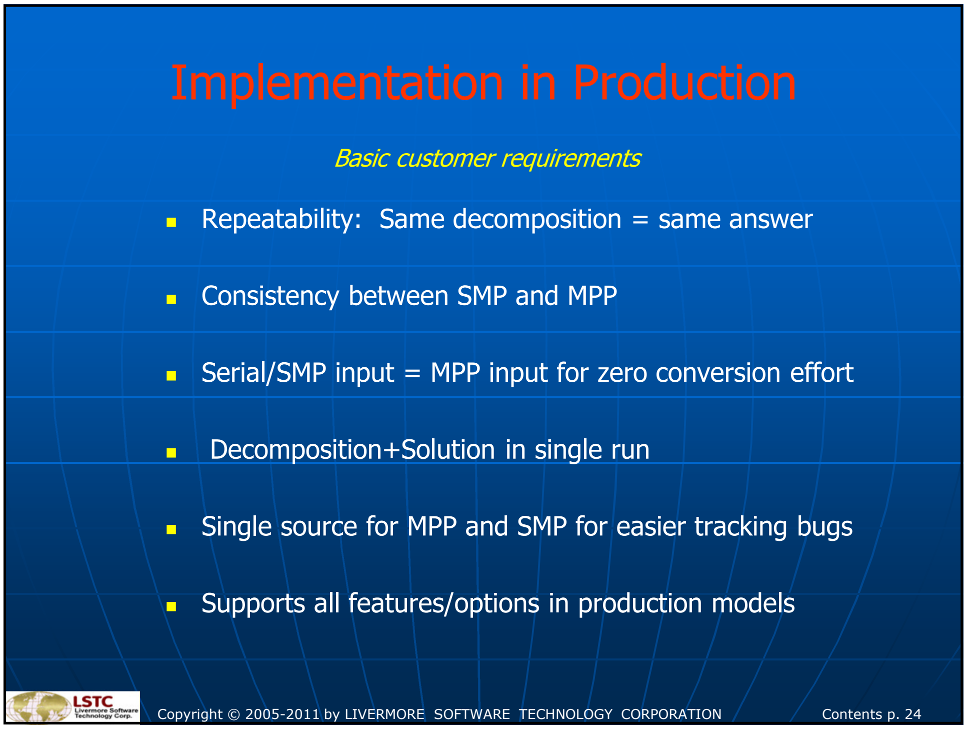
Computing environment



What Drives the MPP Development?

Computing and computer technology

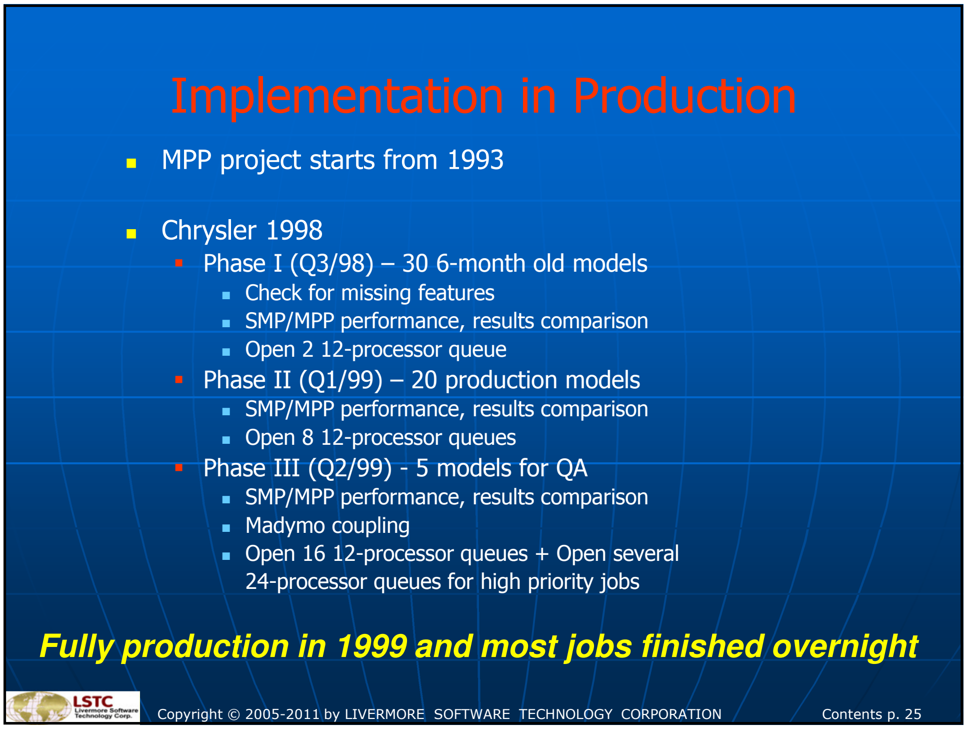




Implementation in Production

Basic customer requirements

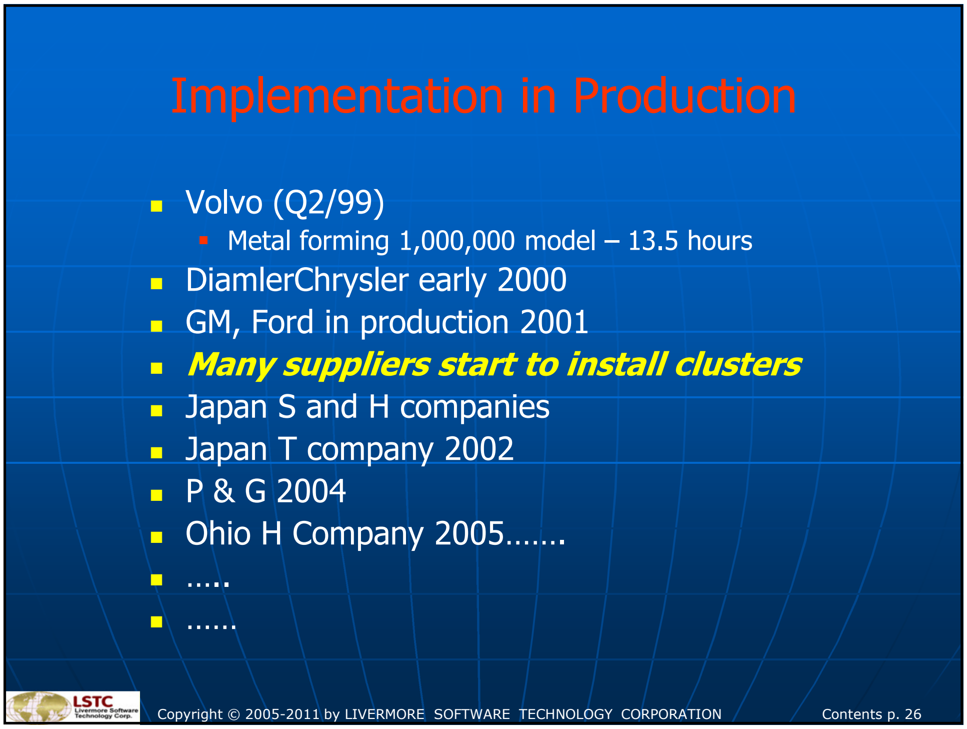
- Repeatability: Same decomposition = same answer
- Consistency between SMP and MPP
- Serial/SMP input = MPP input for zero conversion effort
- Decomposition+Solution in single run
- Single source for MPP and SMP for easier tracking bugs
- Supports all features/options in production models



Implementation in Production

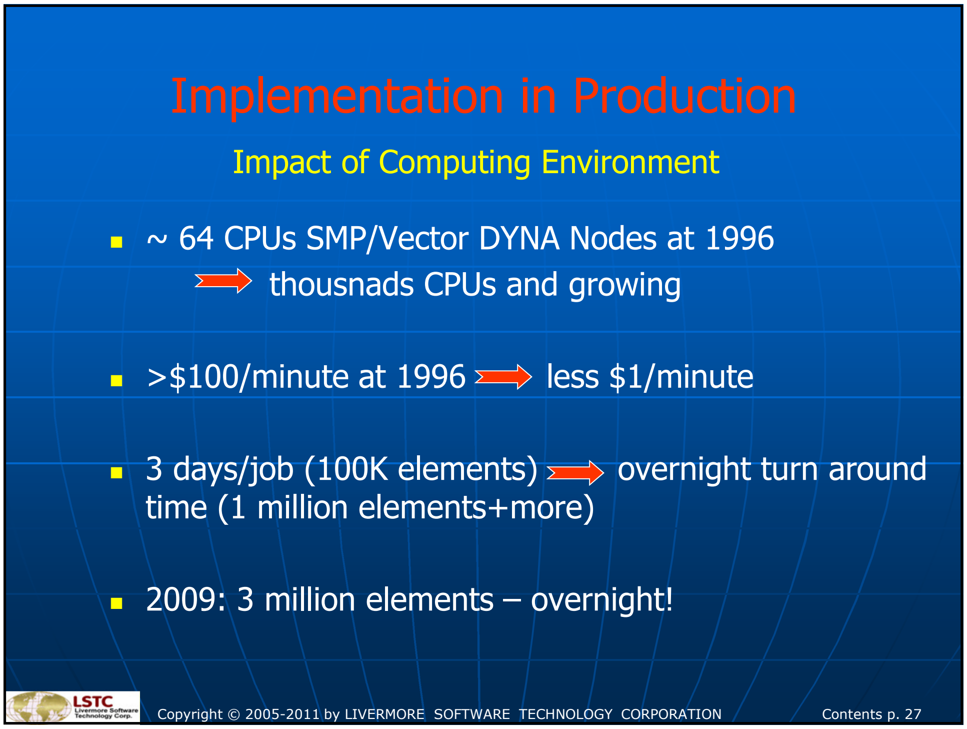
- MPP project starts from 1993
- Chrysler 1998
  - Phase I (Q3/98) – 30 6-month old models
    - Check for missing features
    - SMP/MPP performance, results comparison
    - Open 2 12-processor queue
  - Phase II (Q1/99) – 20 production models
    - SMP/MPP performance, results comparison
    - Open 8 12-processor queues
  - Phase III (Q2/99) - 5 models for QA
    - SMP/MPP performance, results comparison
    - Madymo coupling
    - Open 16 12-processor queues + Open several 24-processor queues for high priority jobs

Fully production in 1999 and most jobs finished overnight



Implementation in Production

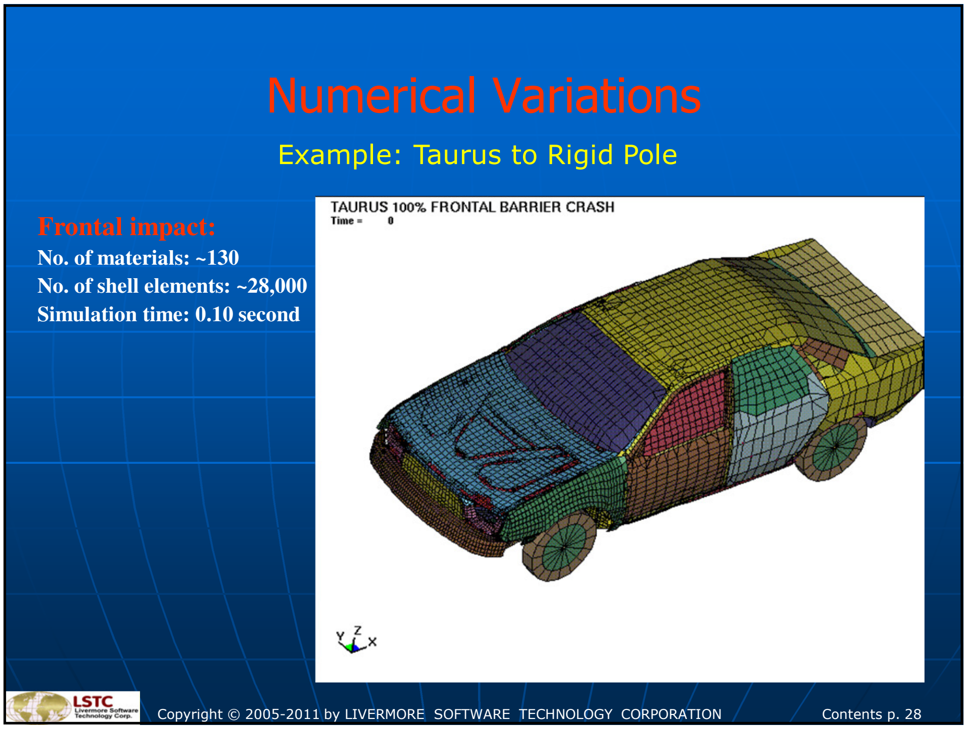
- Volvo (Q2/99)
  - Metal forming 1,000,000 model – 13.5 hours
- DiamlerChrysler early 2000
- GM, Ford in production 2001
- ***Many suppliers start to install clusters***
- Japan S and H companies
- Japan T company 2002
- P & G 2004
- Ohio H Company 2005…….
- …..
- ……



Implementation in Production

Impact of Computing Environment

- ~ 64 CPUs SMP/Vector DYNA Nodes at 1996
  → thousnads CPUs and growing
- >$100/minute at 1996 → less $1/minute
- 3 days/job (100K elements) → overnight turn around time (1 million elements+more)
- 2009: 3 million elements – overnight!



Numerical Variations

Example: Taurus to Rigid Pole

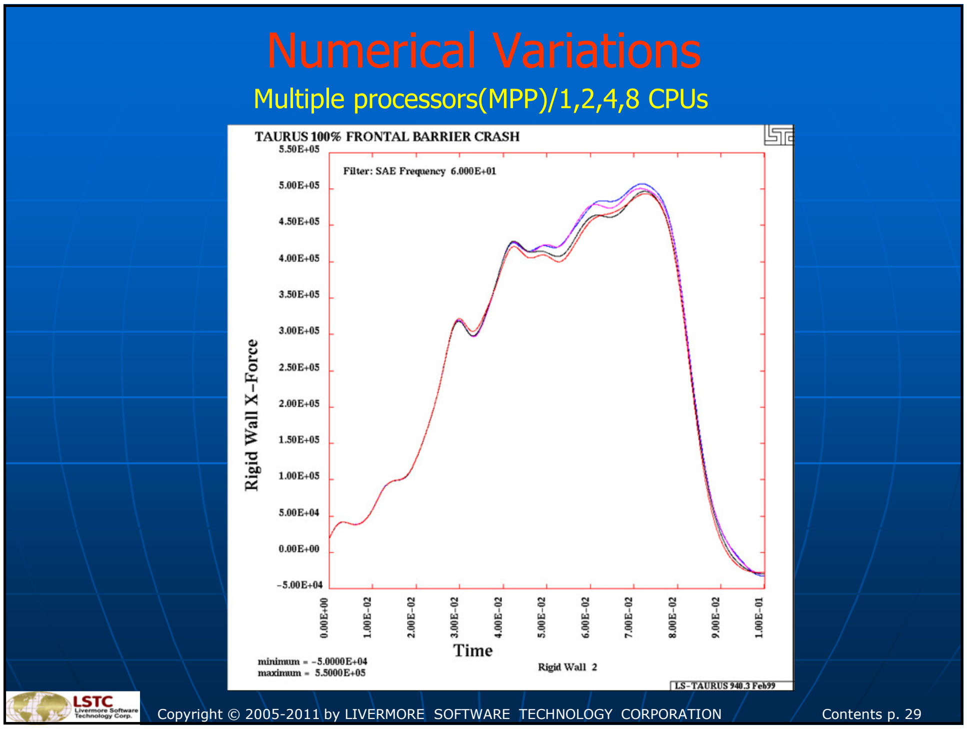
Frontal impact:

No. of materials: ~130

No. of shell elements: ~28,000

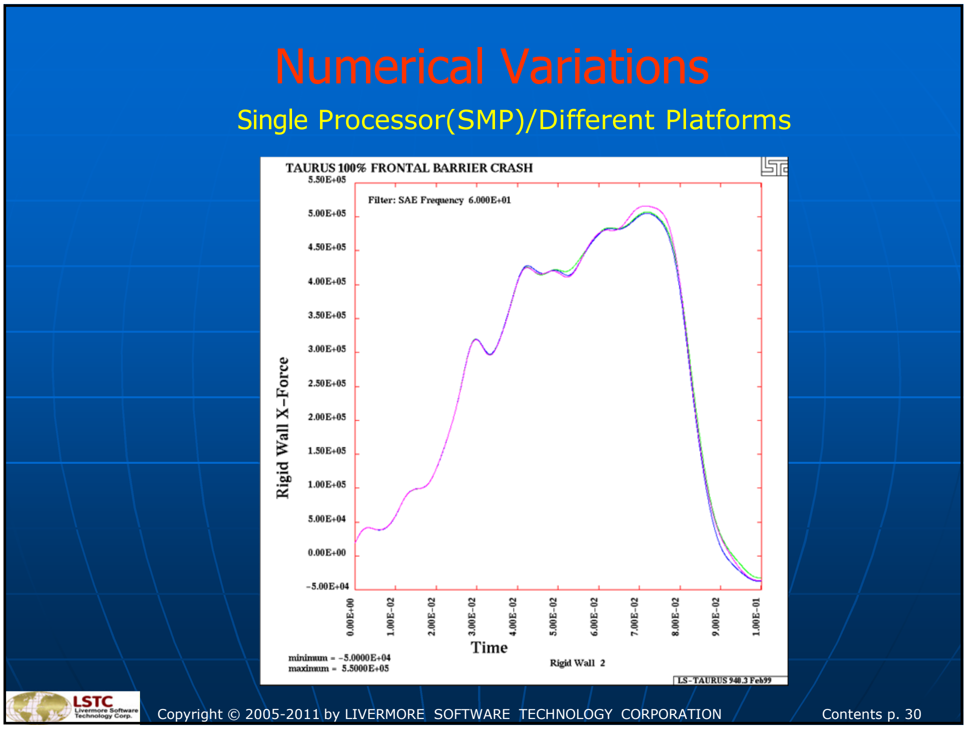
Simulation time: 0.10 second

TAURUS 100% FRONTAL BARRIER CRASH
Time = 0
Y Z X
LSTC Livermore Software Technology Corp.



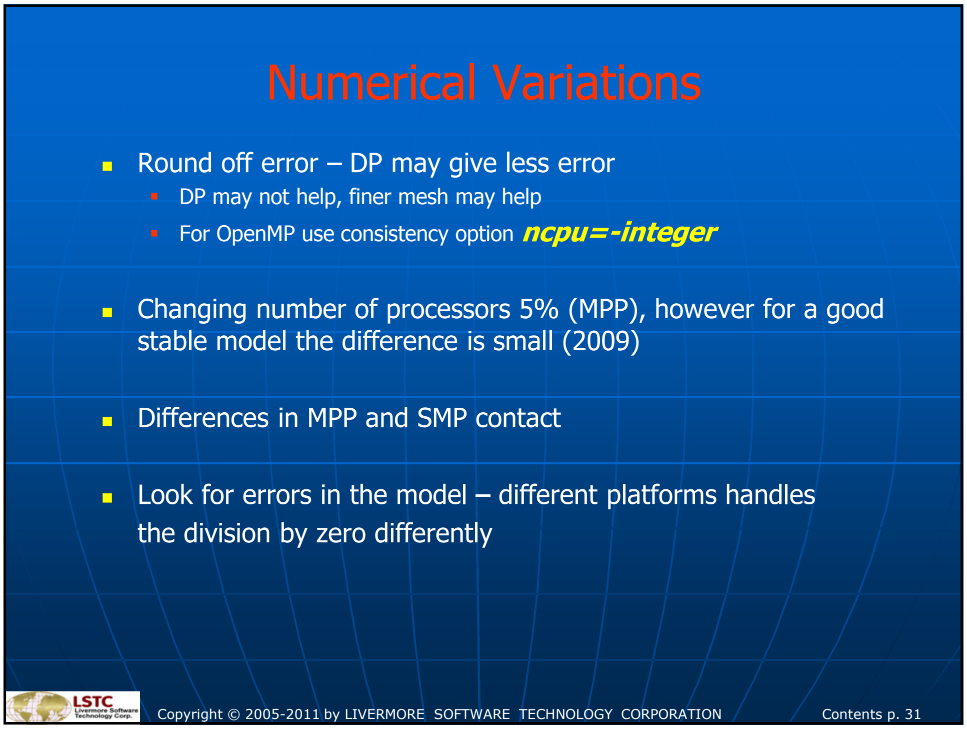
Numerical Variations

Multiple processors(MPP)/1,2,4,8 CPUs



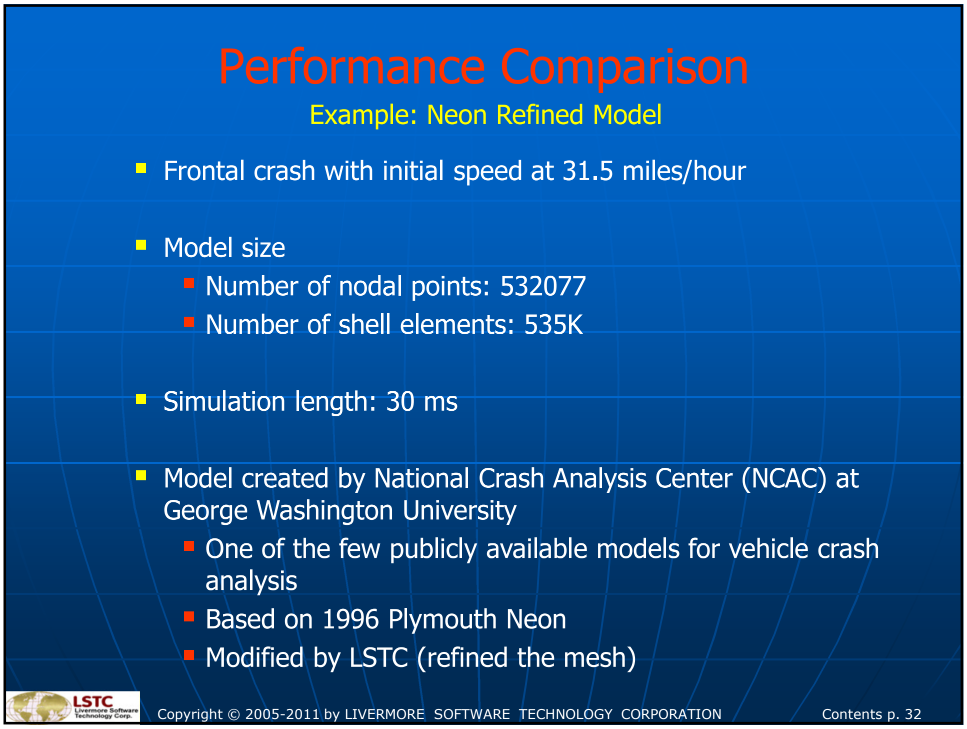
Numerical Variations

Single Processor(SMP)/Different Platforms



Numerical Variations

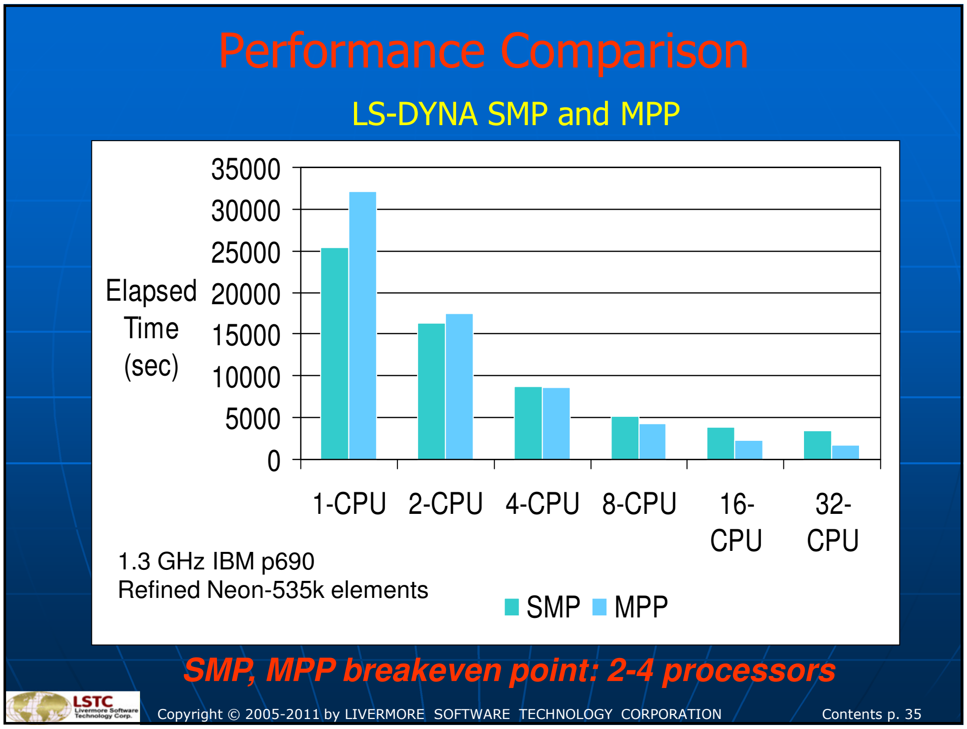
- Round off error – DP may give less error
  - DP may not help, finer mesh may help
  - For OpenMP use consistency option ***ncpu=-integer***
- Changing number of processors 5% (MPP), however for a good stable model the difference is small (2009)
- Differences in MPP and SMP contact
- Look for errors in the model – different platforms handles the division by zero differently



Performance Comparison

Example: Neon Refined Model

- Frontal crash with initial speed at 31.5 miles/hour
- Model size
  - Number of nodal points: 532077
  - Number of shell elements: 535K
- Simulation length: 30 ms
- Model created by National Crash Analysis Center (NCAC) at George Washington University
  - One of the few publicly available models for vehicle crash analysis
  - Based on 1996 Plymouth Neon
  - Modified by LSTC (refined the mesh)

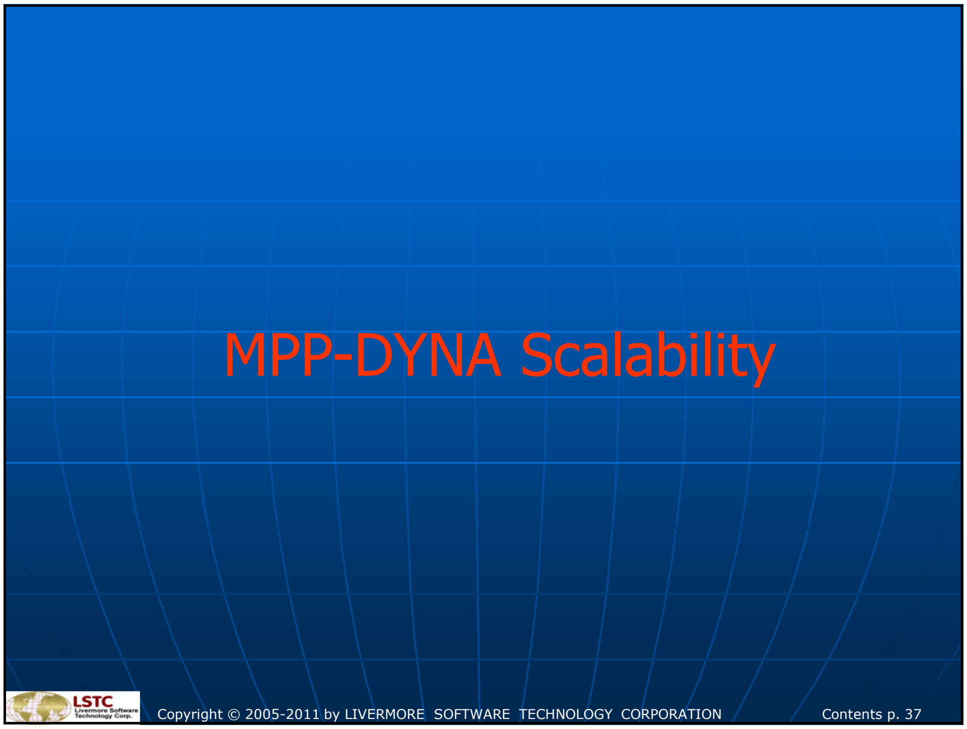


Performance Comparison

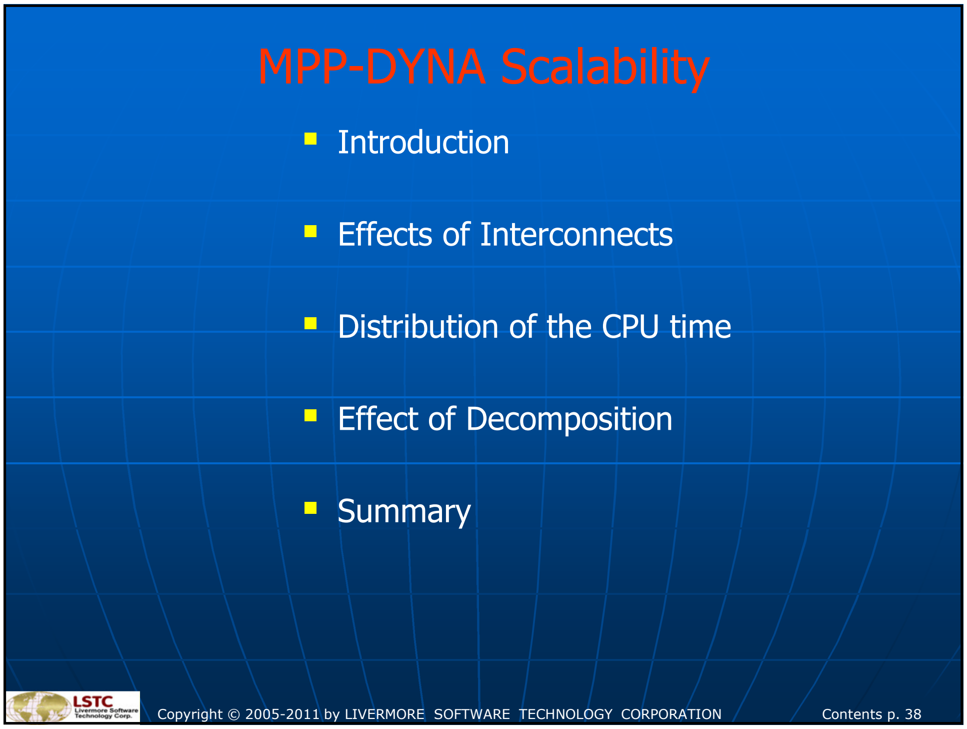
LS-DYNA SMP and MPP

1.3 GHz IBM p690
Refined Neon-535k elements

SMP, MPP breakeven point: 2-4 processors

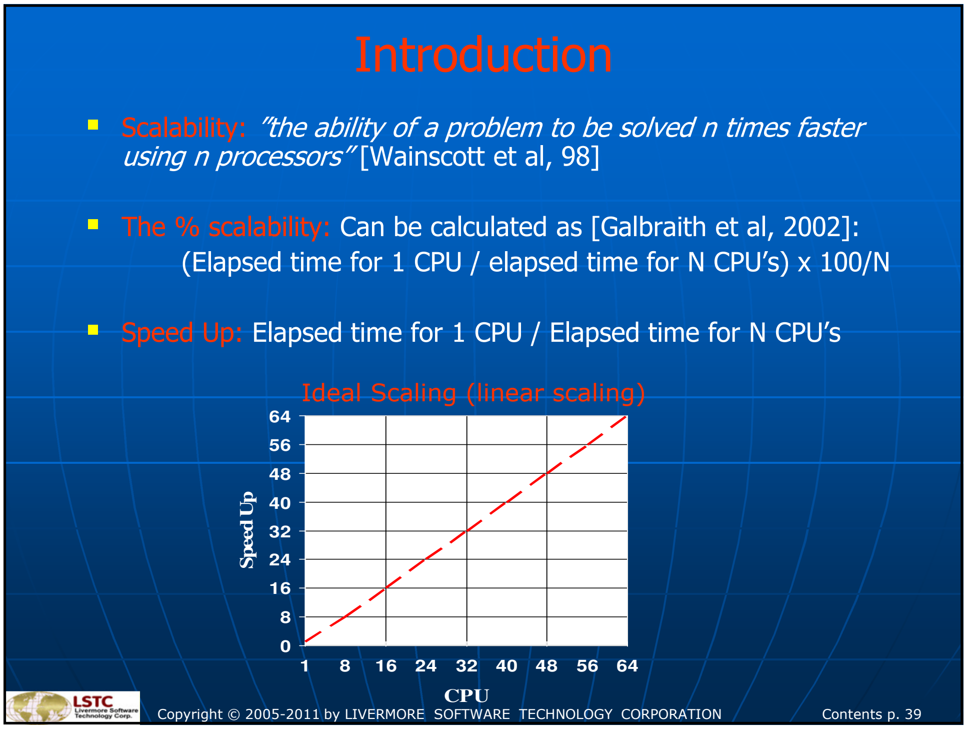


MPP-DYNA Scalability



MPP-DYNA Scalability

- Introduction
- Effects of Interconnects
- Distribution of the CPU time
- Effect of Decomposition
- Summary



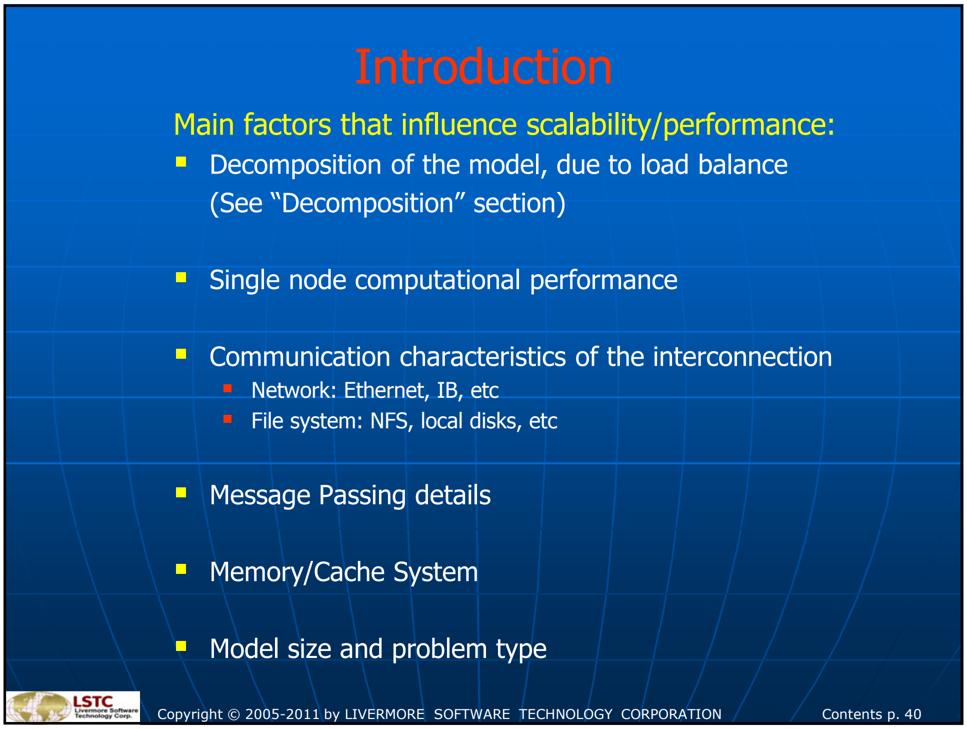
Introduction

- Scalability: *"the ability of a problem to be solved n times faster using n processors"* [Wainscott et al, 98]

- The % scalability: Can be calculated as [Galbraith et al, 2002]:
  (Elapsed time for 1 CPU / elapsed time for N CPU's) x 100/N

- Speed Up: Elapsed time for 1 CPU / Elapsed time for N CPU's

Ideal Scaling (linear scaling)



Introduction

Main factors that influence scalability/performance:

- Decomposition of the model, due to load balance
  (See "Decomposition" section)

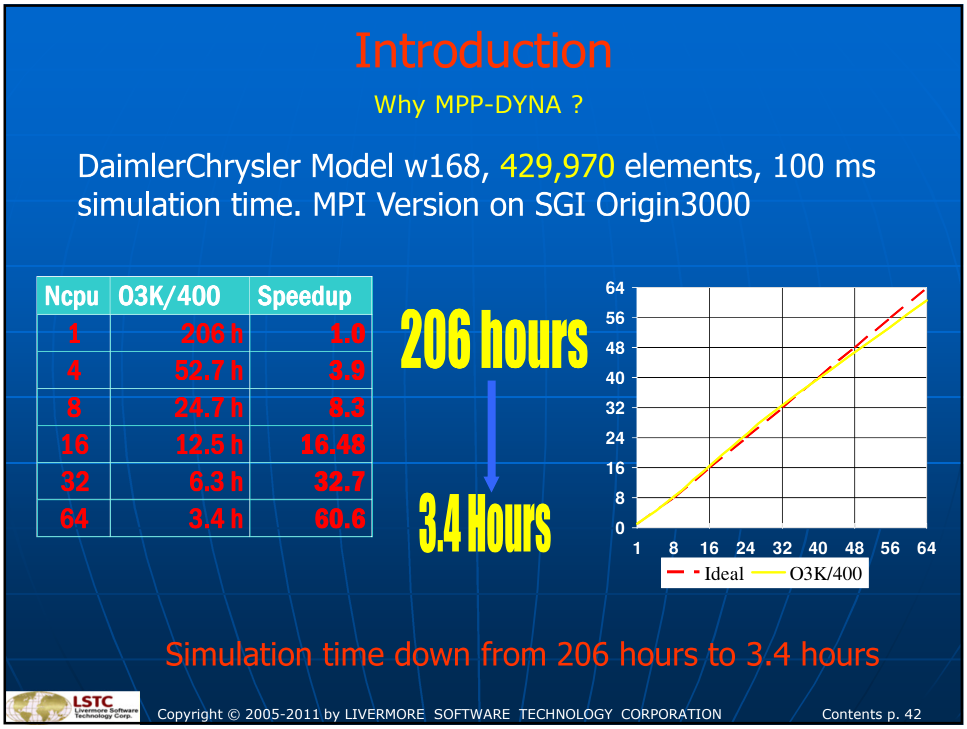
- Single node computational performance

- Communication characteristics of the interconnection
  - Network: Ethernet, IB, etc
  - File system: NFS, local disks, etc

- Message Passing details

- Memory/Cache System

- Model size and problem type



Introduction

Why MPP-DYNA ?

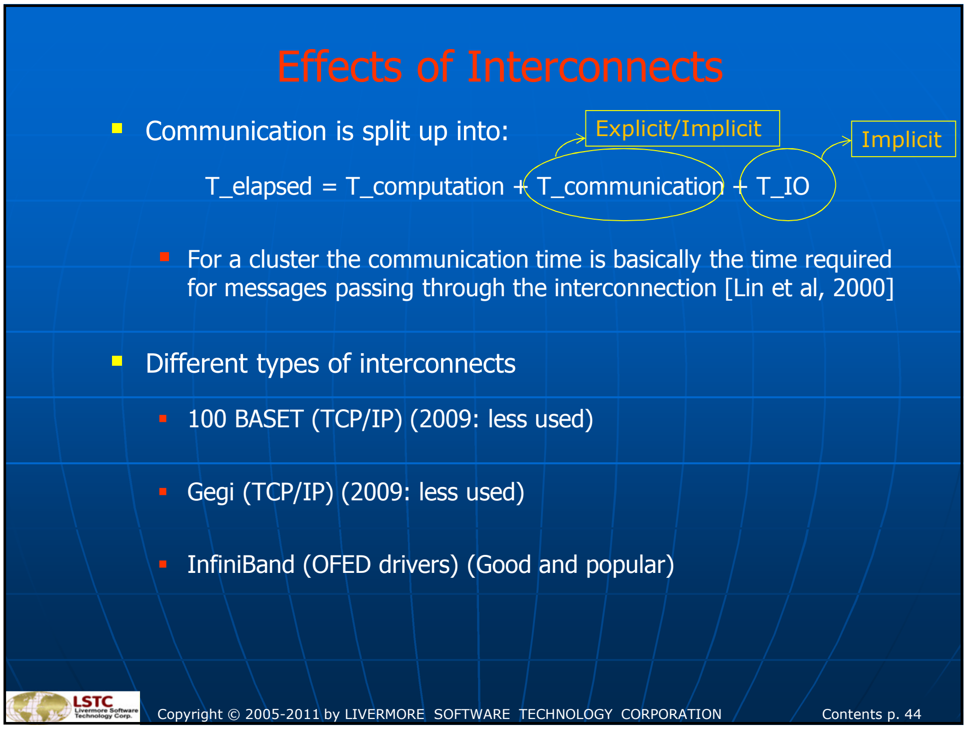
DaimlerChrysler Model w168, 429,970 elements, 100 ms simulation time. MPI Version on SGI Origin3000

Ncpu	O3K/400	Speedup
1	206 h	1.0
4	52.7 h	3.9
8	24.7 h	8.3
16	12.5 h	16.48
32	6.3 h	32.7
64	3.4 h	60.6

206 hours

3.4 Hours

Simulation time down from 206 hours to 3.4 hours



Effects of Interconnects

- Communication is split up into:

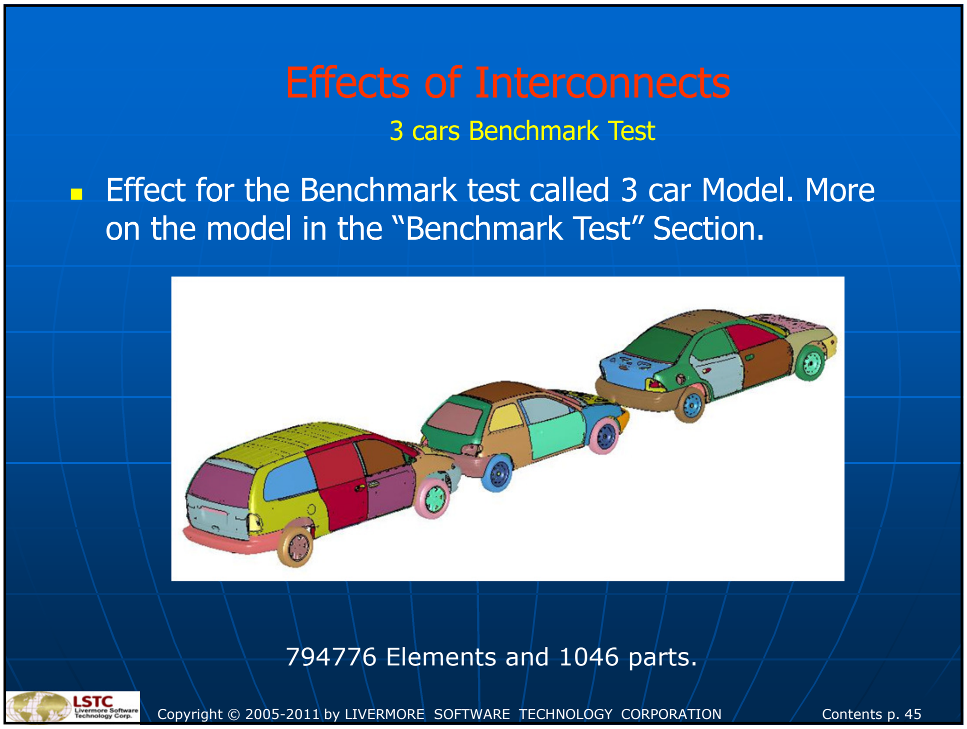
  T_elapsed = T_computation + T_communication + T_IO

  Explicit/Implicit

  Implicit

  - For a cluster the communication time is basically the time required for messages passing through the interconnection [Lin et al, 2000]

- Different types of interconnects
  - 100 BASET (TCP/IP) (2009: less used)
  - Gegi (TCP/IP) (2009: less used)
  - InfiniBand (OFED drivers) (Good and popular)

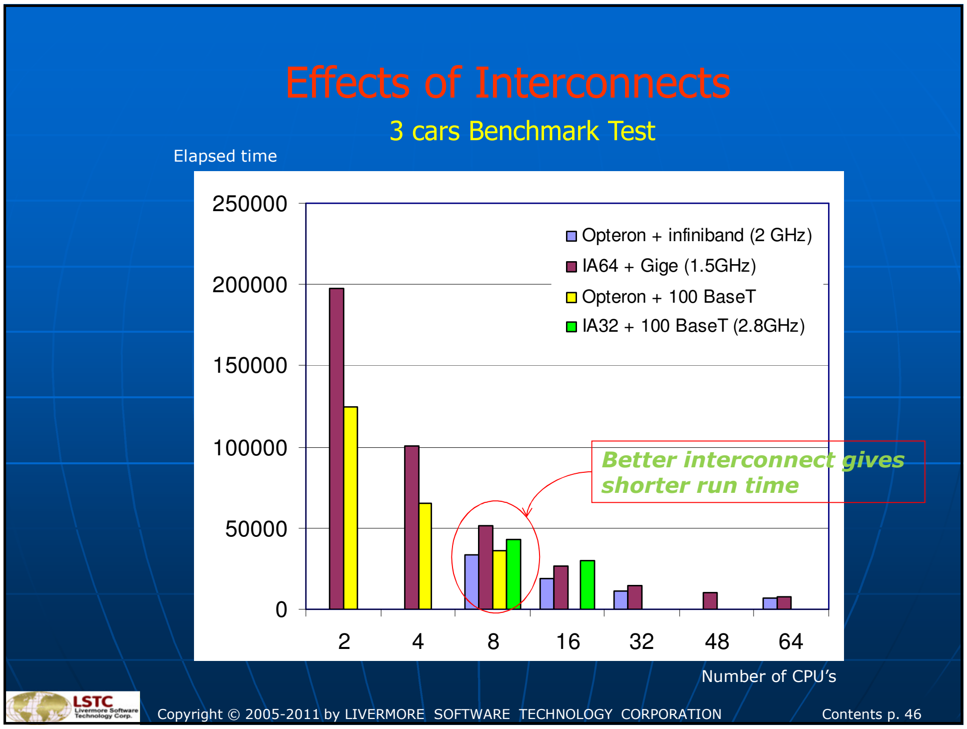


Effects of Interconnects

3 cars Benchmark Test

- Effect for the Benchmark test called 3 car Model. More on the model in the “Benchmark Test” Section.

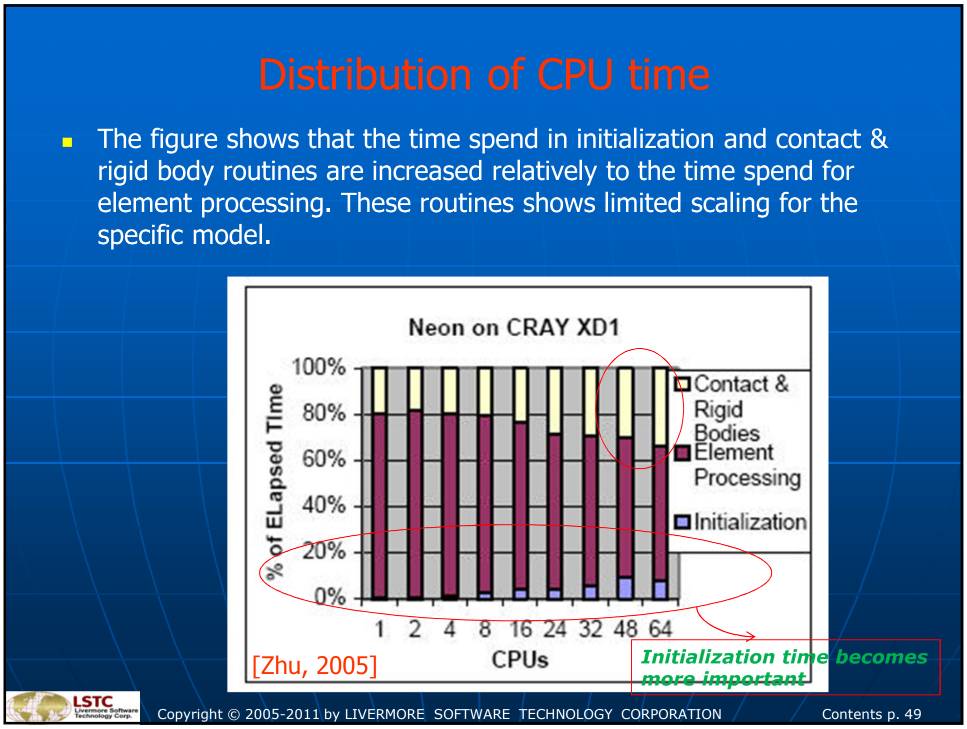
794776 Elements and 1046 parts.



Effects of Interconnects

3 cars Benchmark Test

Better interconnect gives shorter run time



Distribution of CPU time

- The figure shows that the time spend in initialization and contact & rigid body routines are increased relatively to the time spend for element processing. These routines shows limited scaling for the specific model.

[Zhu, 2005]

Initialization time becomes more important

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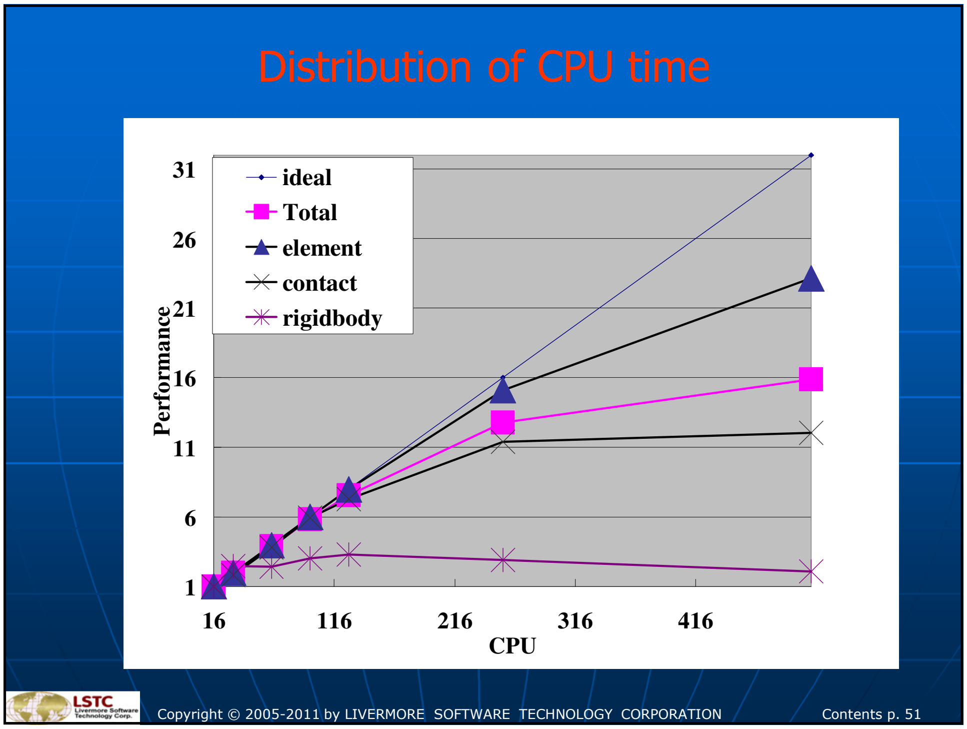

Distribution of CPU time

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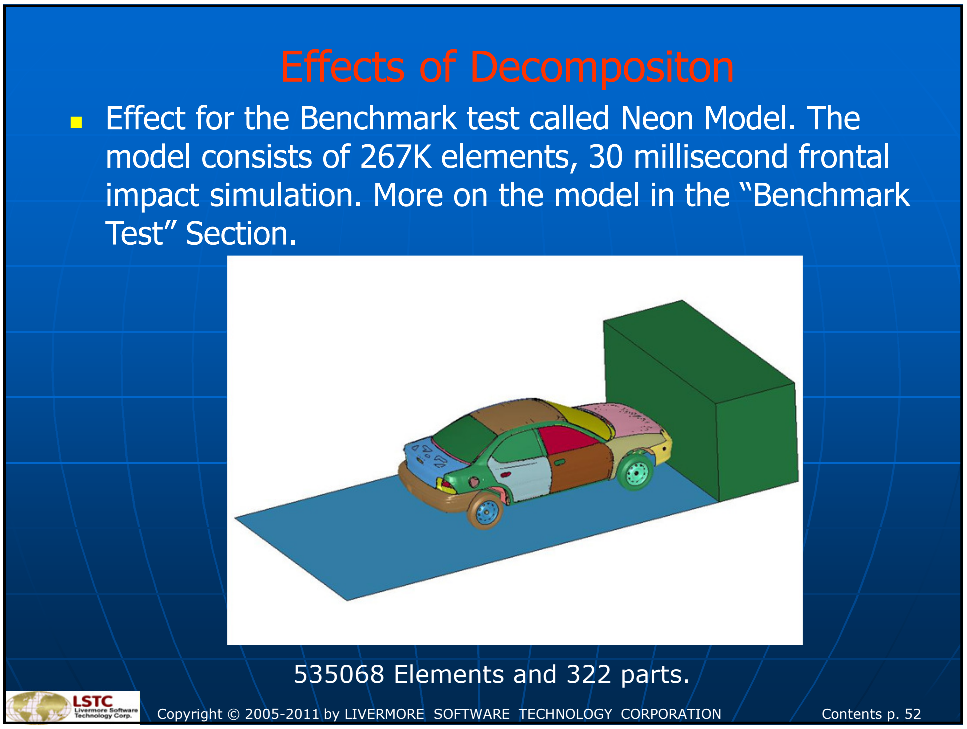
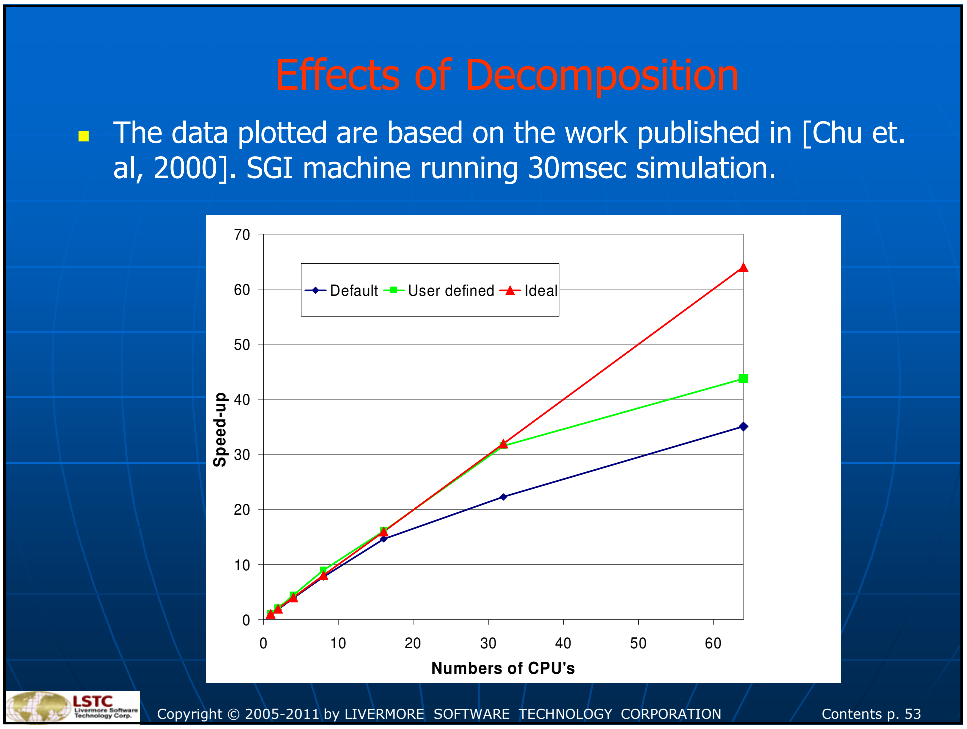


Effects of Decompositon

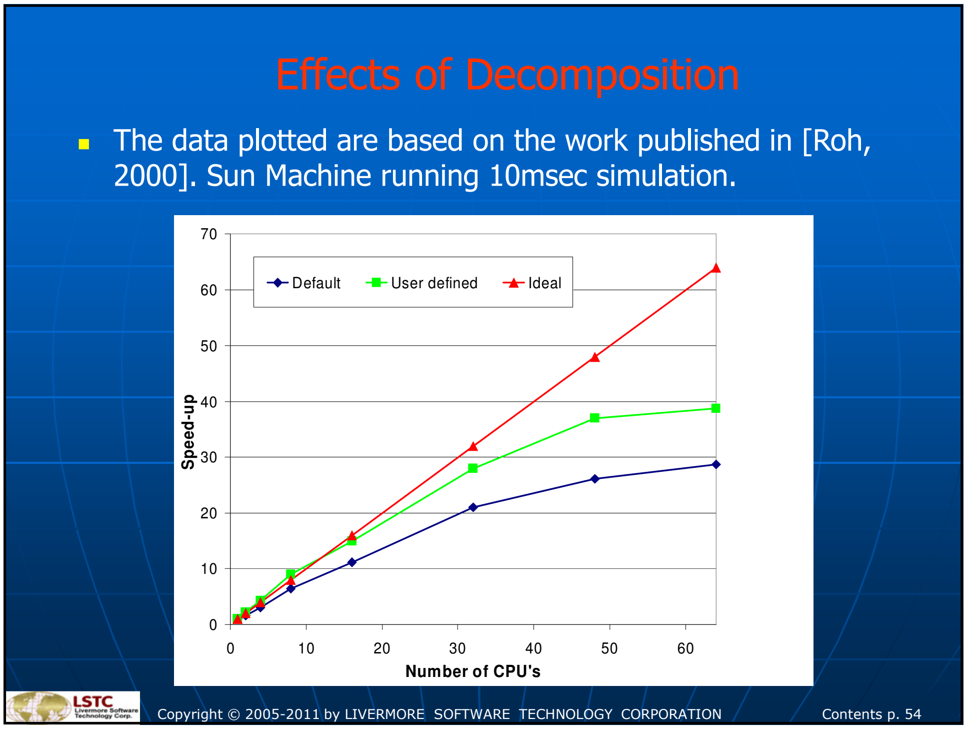
- Effect for the Benchmark test called Neon Model. The model consists of 267K elements, 30 millisecond frontal impact simulation. More on the model in the “Benchmark Test” Section.

535068 Elements and 322 parts.



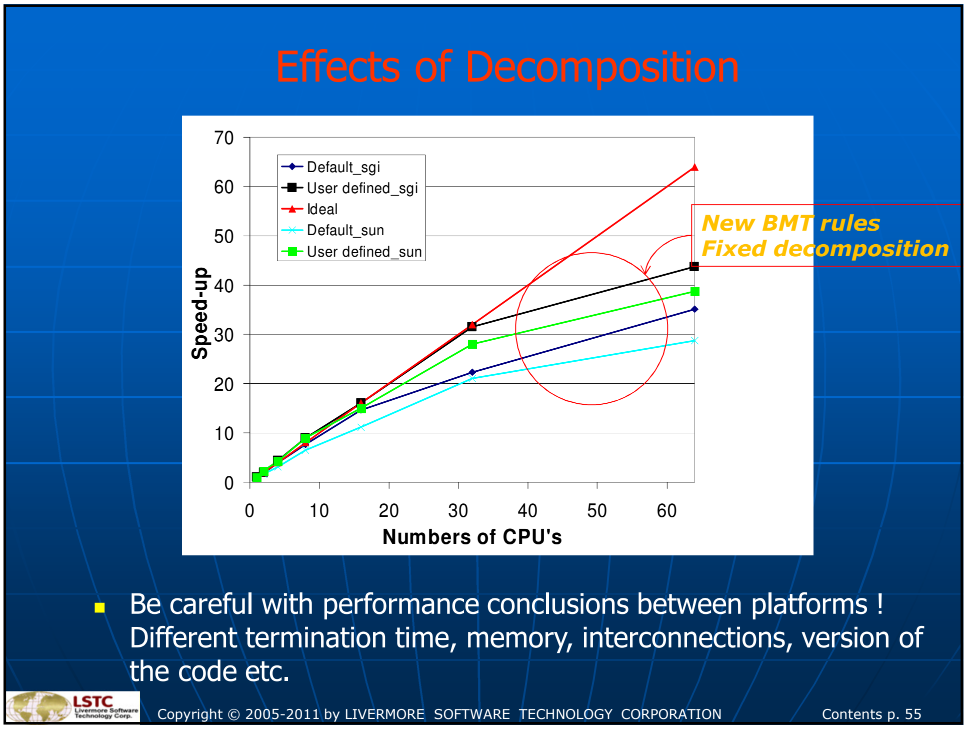
Effects of Decomposition

- The data plotted are based on the work published in [Chu et. al, 2000]. SGI machine running 30msec simulation.



Effects of Decomposition

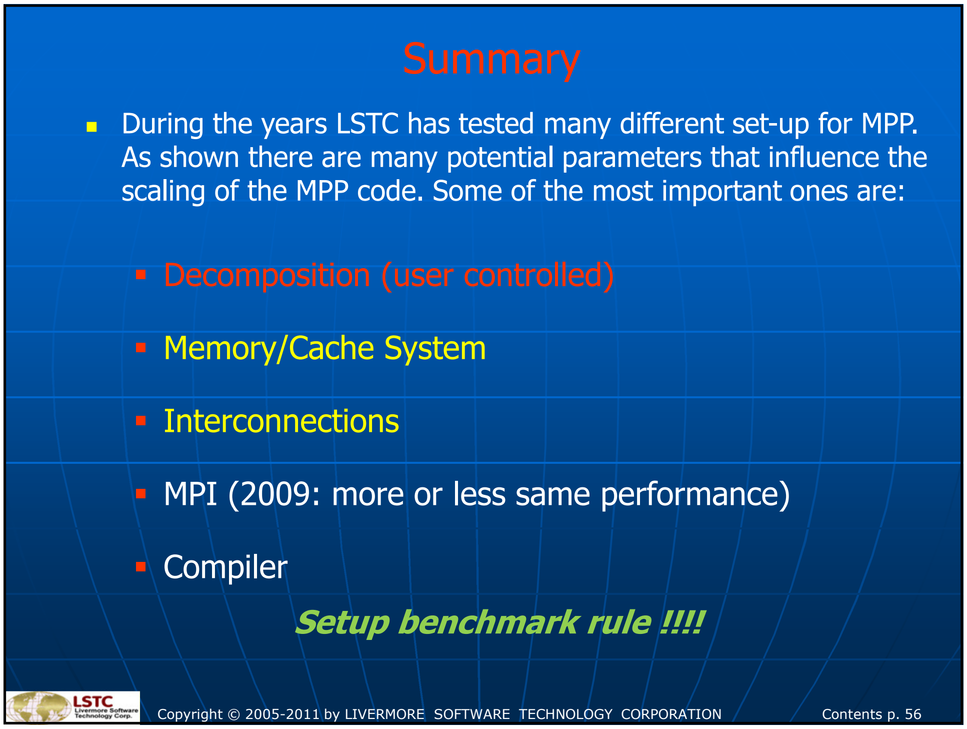
- The data plotted are based on the work published in [Roh, 2000]. Sun Machine running 10msec simulation.



Effects of Decomposition

New BMT rules
Fixed decomposition

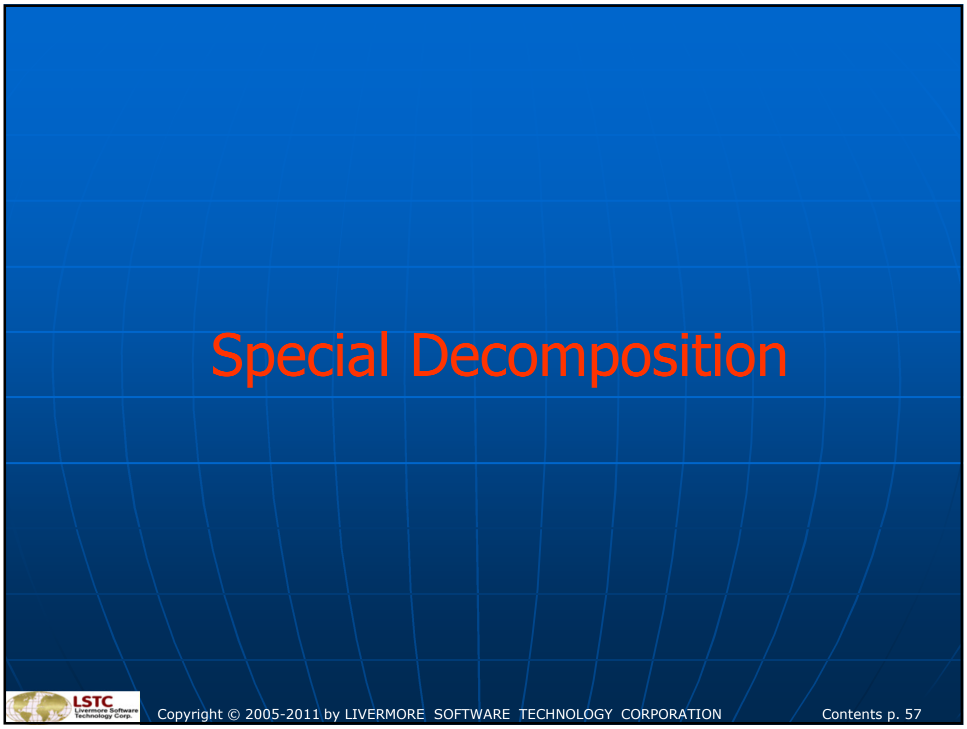
- Be careful with performance conclusions between platforms ! Different termination time, memory, interconnections, version of the code etc.



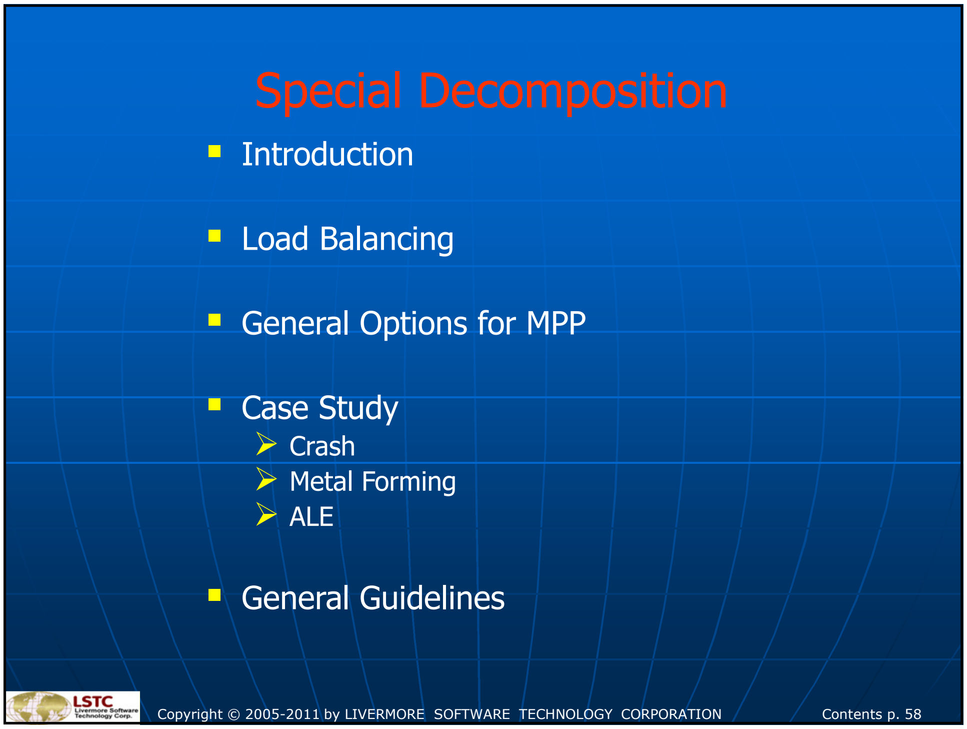
Summary

- During the years LSTC has tested many different set-up for MPP. As shown there are many potential parameters that influence the scaling of the MPP code. Some of the most important ones are:
  - Decomposition (user controlled)
  - Memory/Cache System
  - Interconnections
  - MPI (2009: more or less same performance)
  - Compiler

Setup benchmark rule !!!!

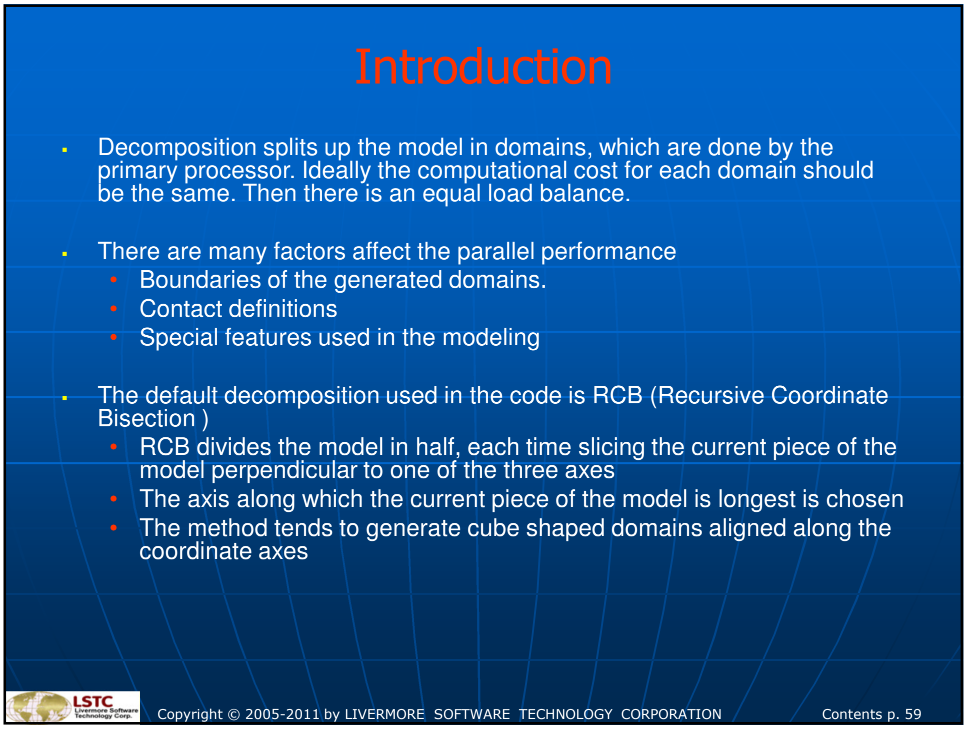


Special Decomposition



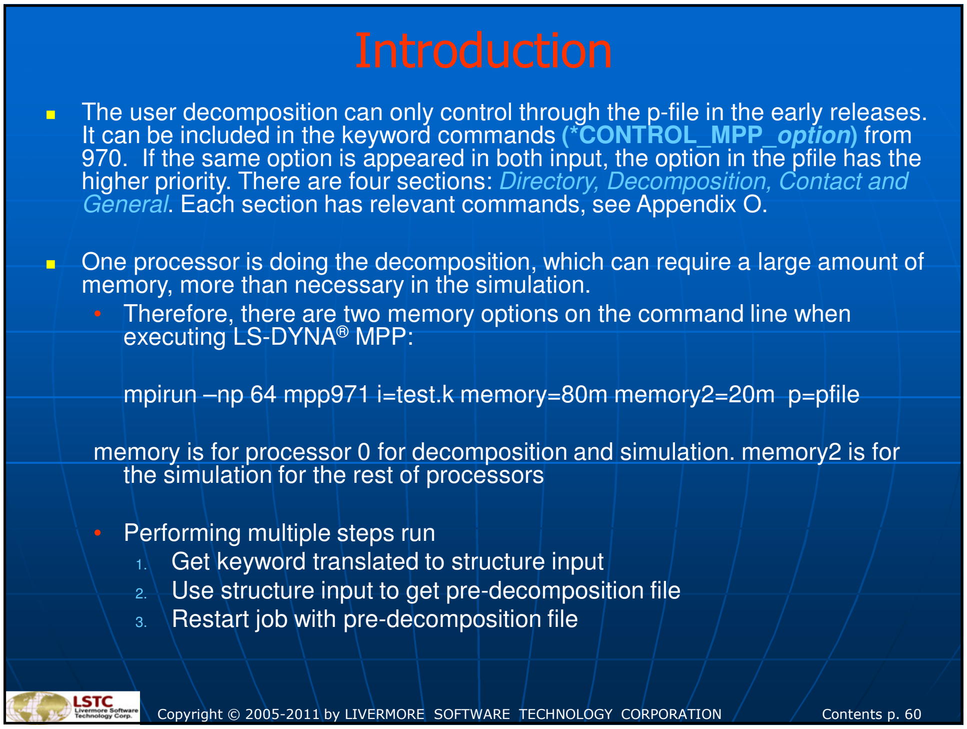
Special Decomposition

- Introduction
- Load Balancing
- General Options for MPP
- Case Study
  - Crash
  - Metal Forming
  - ALE
- General Guidelines



Introduction

- Decomposition splits up the model in domains, which are done by the primary processor. Ideally the computational cost for each domain should be the same. Then there is an equal load balance.
- There are many factors affect the parallel performance
  - Boundaries of the generated domains.
  - Contact definitions
  - Special features used in the modeling
- The default decomposition used in the code is RCB (Recursive Coordinate Bisection)
  - RCB divides the model in half, each time slicing the current piece of the model perpendicular to one of the three axes
  - The axis along which the current piece of the model is longest is chosen
  - The method tends to generate cube shaped domains aligned along the coordinate axes



Introduction

- The user decomposition can only control through the p-file in the early releases. It can be included in the keyword commands **(*CONTROL_MPP_*option*)** from 970. If the same option is appeared in both input, the option in the pfile has the higher priority. There are four sections: *Directory, Decomposition, Contact and General*. Each section has relevant commands, see Appendix O.

- One processor is doing the decomposition, which can require a large amount of memory, more than necessary in the simulation.
  - Therefore, there are two memory options on the command line when executing LS-DYNA® MPP:

    mpirun –np 64 mpp971 i=test.k memory=80m memory2=20m p=pfile

  memory is for processor 0 for decomposition and simulation. memory2 is for the simulation for the rest of processors

  - Performing multiple steps run
    1. Get keyword translated to structure input
    2. Use structure input to get pre-decomposition file
    3. Restart job with pre-decomposition file

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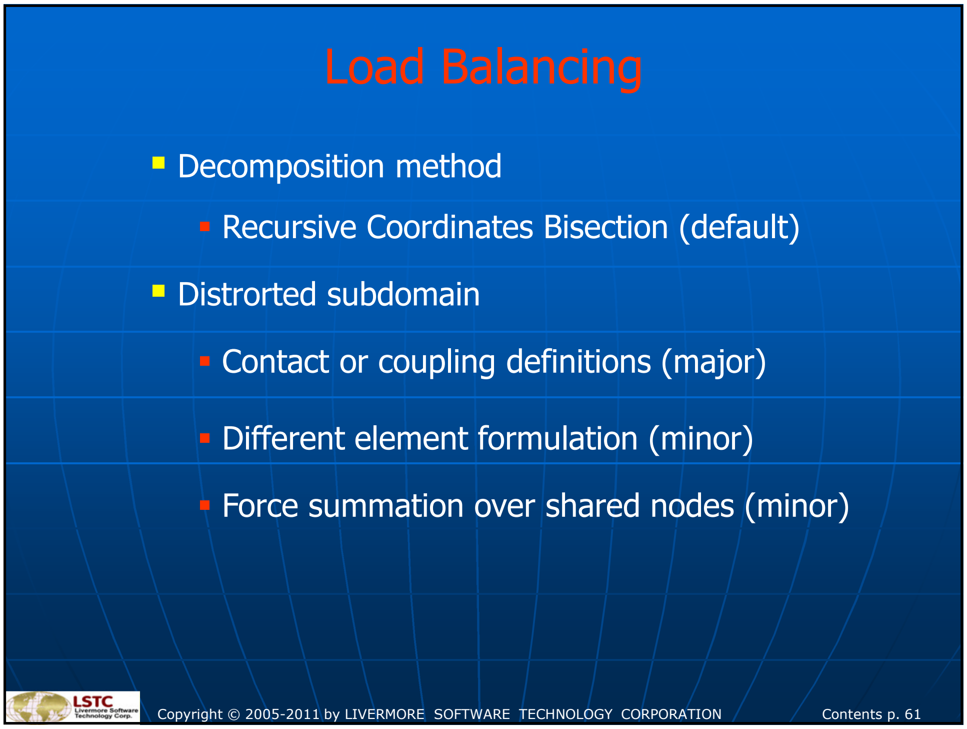

Load Balancing

- Decomposition method
  - Recursive Coordinates Bisection (default)
- Distrorted subdomain
  - Contact or coupling definitions (major)
  - Different element formulation (minor)
  - Force summation over shared nodes (minor)

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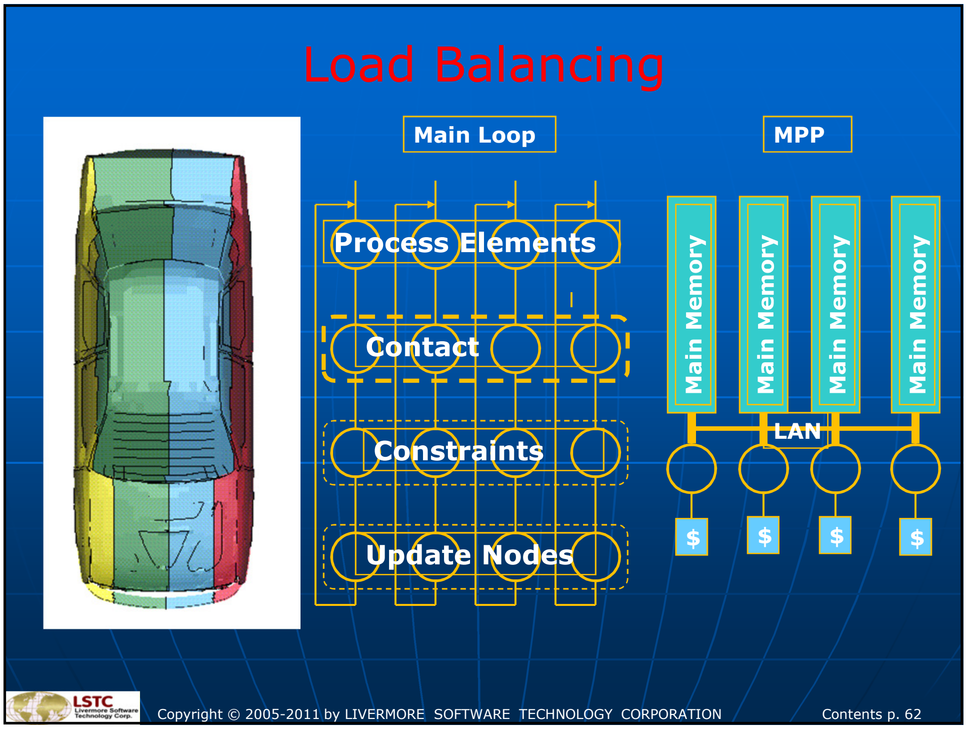
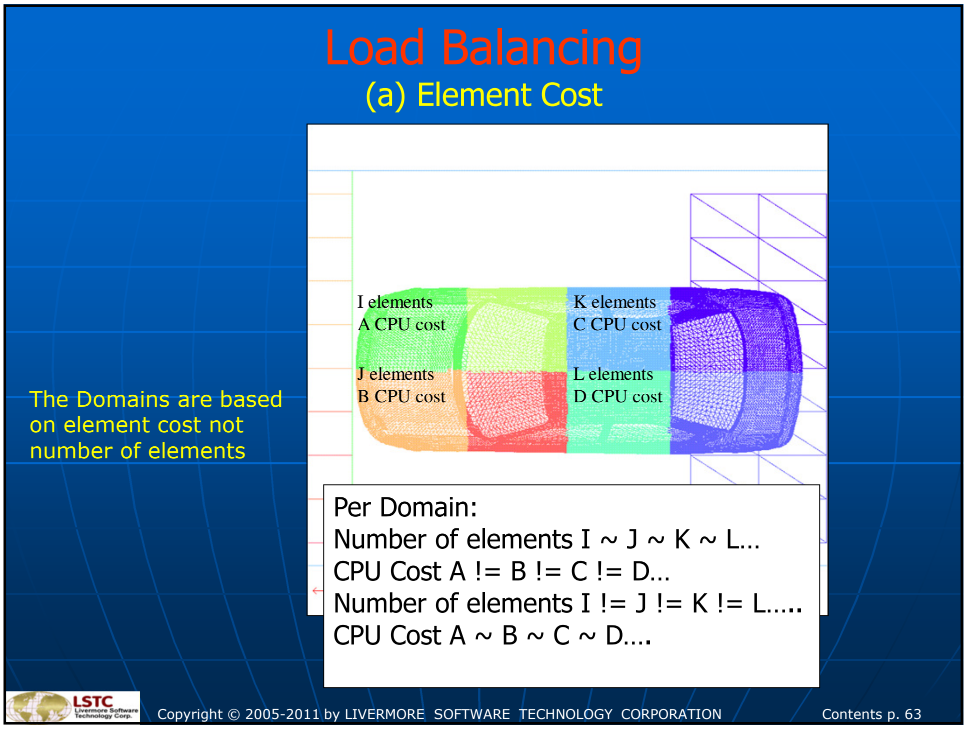


Load Balancing

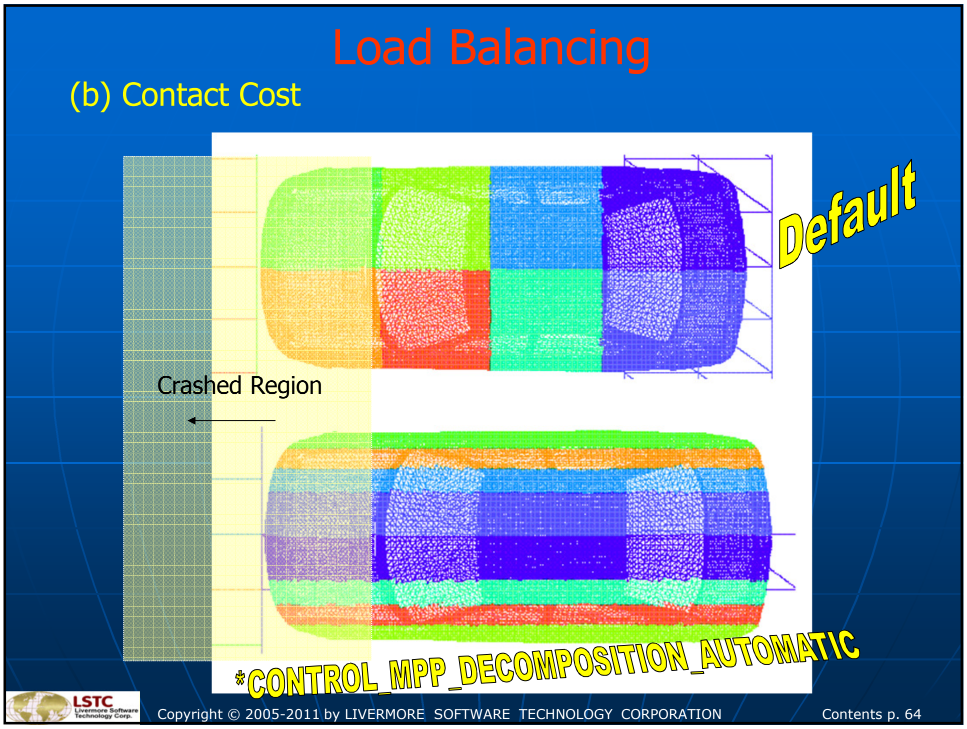


Load Balancing

(a) Element Cost

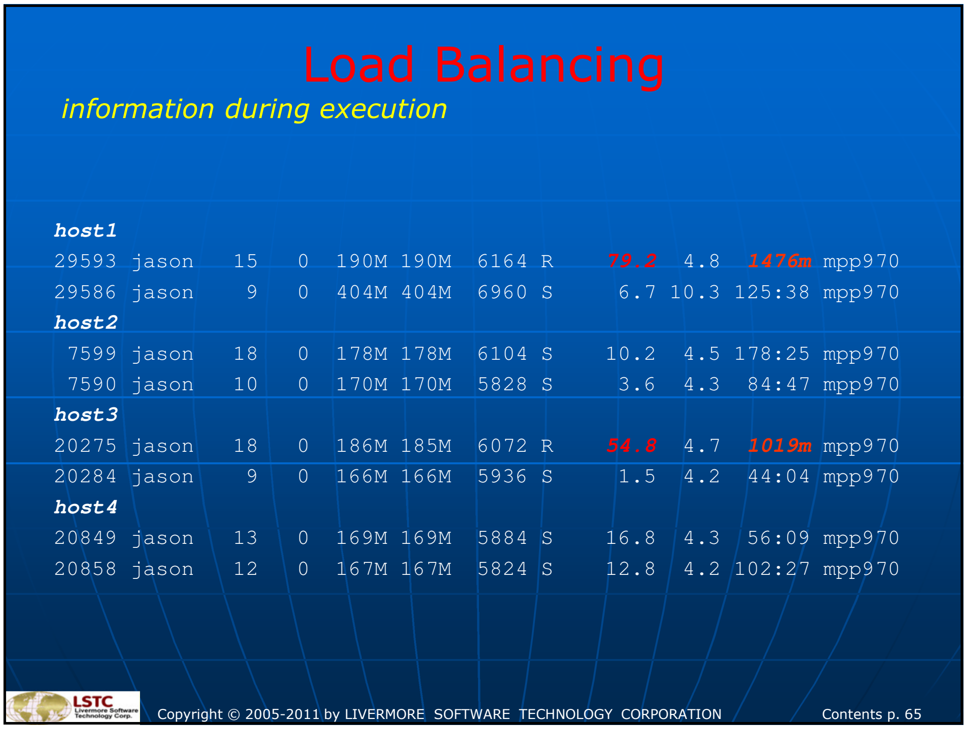
The Domains are based on element cost not number of elements

Per Domain:
Number of elements I ~ J ~ K ~ L…
CPU Cost A != B != C != D…
Number of elements I != J != K != L…..
CPU Cost A ~ B ~ C ~ D….



Load Balancing

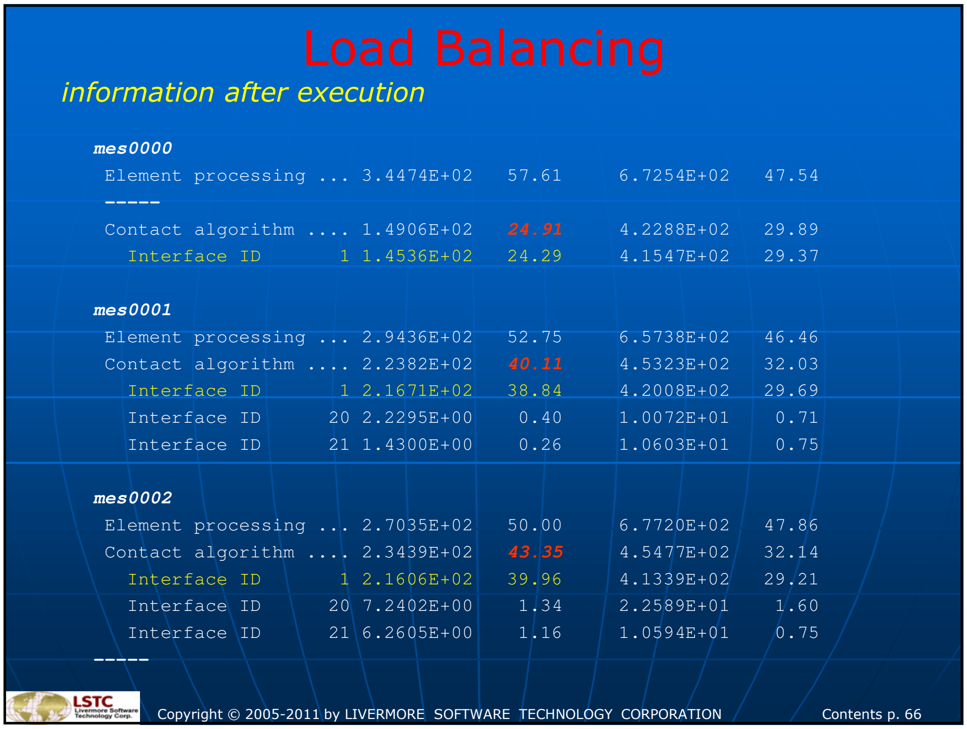
(b) Contact Cost



Load Balancing

information during execution

```
host1
29593 jason    15    0   190M 190M   6164 R      79.2   4.8   1476m mpp970
29586 jason     9    0   404M 404M   6960 S       6.7  10.3 125:38 mpp970
host2
 7599 jason    18    0   178M 178M   6104 S      10.2   4.5 178:25 mpp970
 7590 jason    10    0   170M 170M   5828 S       3.6   4.3   84:47 mpp970
host3
20275 jason    18    0   186M 185M   6072 R      54.8   4.7   1019m mpp970
20284 jason     9    0   166M 166M   5936 S       1.5   4.2   44:04 mpp970
host4
20849 jason    13    0   169M 169M   5884 S      16.8   4.3   56:09 mpp970
20858 jason    12    0   167M 167M   5824 S      12.8   4.2 102:27 mpp970
```



Load Balancing

information after execution

```
mes0000
 Element processing ... 3.4474E+02   57.61    6.7254E+02   47.54
 -----
 Contact algorithm .... 1.4906E+02   24.91    4.2288E+02   29.89
   Interface ID       1 1.4536E+02   24.29    4.1547E+02   29.37

mes0001
 Element processing ... 2.9436E+02   52.75    6.5738E+02   46.46
 Contact algorithm .... 2.2382E+02   40.11    4.5323E+02   32.03
   Interface ID       1 2.1671E+02   38.84    4.2008E+02   29.69
   Interface ID      20 2.2295E+00    0.40    1.0072E+01    0.71
   Interface ID      21 1.4300E+00    0.26    1.0603E+01    0.75

mes0002
 Element processing ... 2.7035E+02   50.00    6.7720E+02   47.86
 Contact algorithm .... 2.3439E+02   43.35    4.5477E+02   32.14
   Interface ID       1 2.1606E+02   39.96    4.1339E+02   29.21
   Interface ID      20 7.2402E+00    1.34    2.2589E+01    1.60
   Interface ID      21 6.2605E+00    1.16    1.0594E+01    0.75
-----
```

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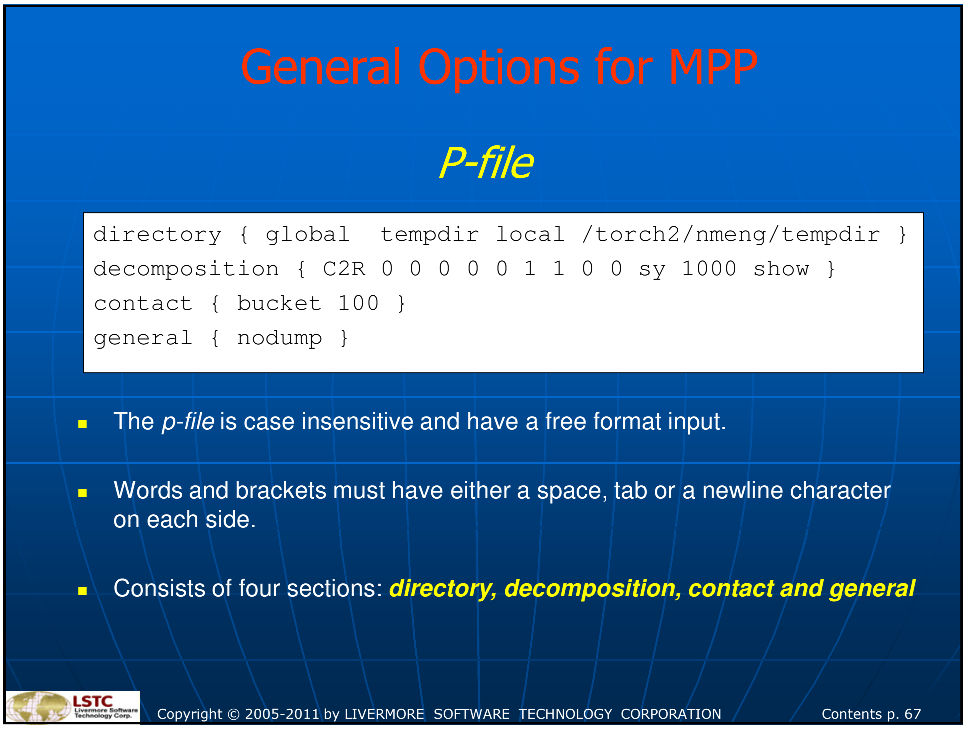

General Options for MPP

P-file

```
directory { global  tempdir local /torch2/nmeng/tempdir }
decomposition { C2R 0 0 0 0 0 1 1 0 0 sy 1000 show }
contact { bucket 100 }
general { nodump }
```

- The *p-file* is case insensitive and have a free format input.
- Words and brackets must have either a space, tab or a newline character on each side.
- Consists of four sections: ***directory, decomposition, contact and general***

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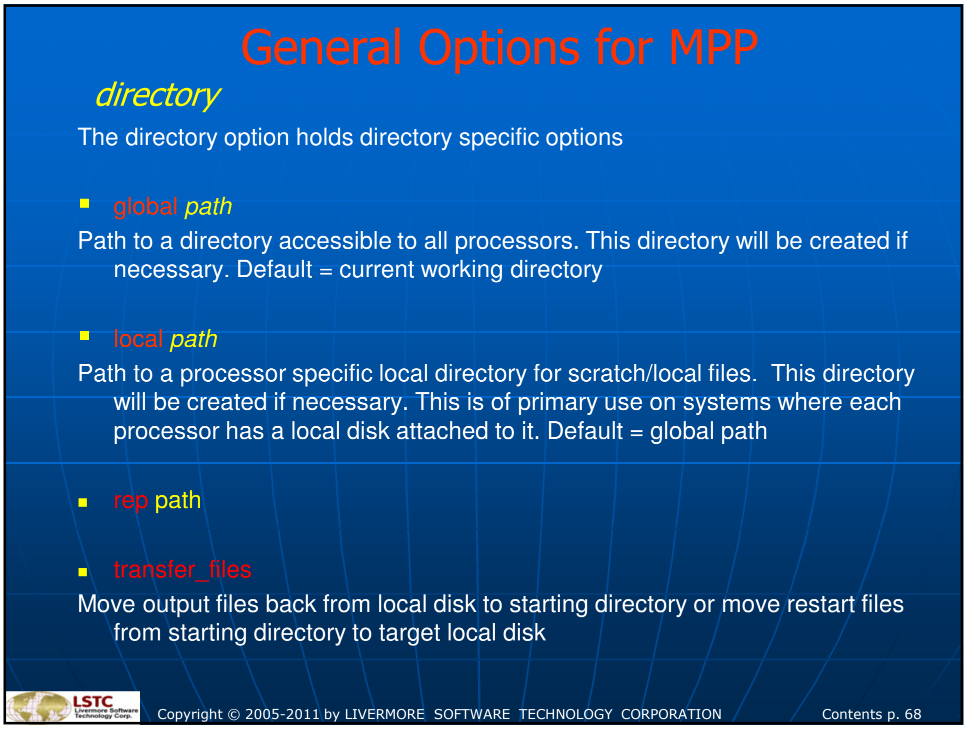


General Options for MPP

directory

The directory option holds directory specific options

- global *path*

Path to a directory accessible to all processors. This directory will be created if necessary. Default = current working directory

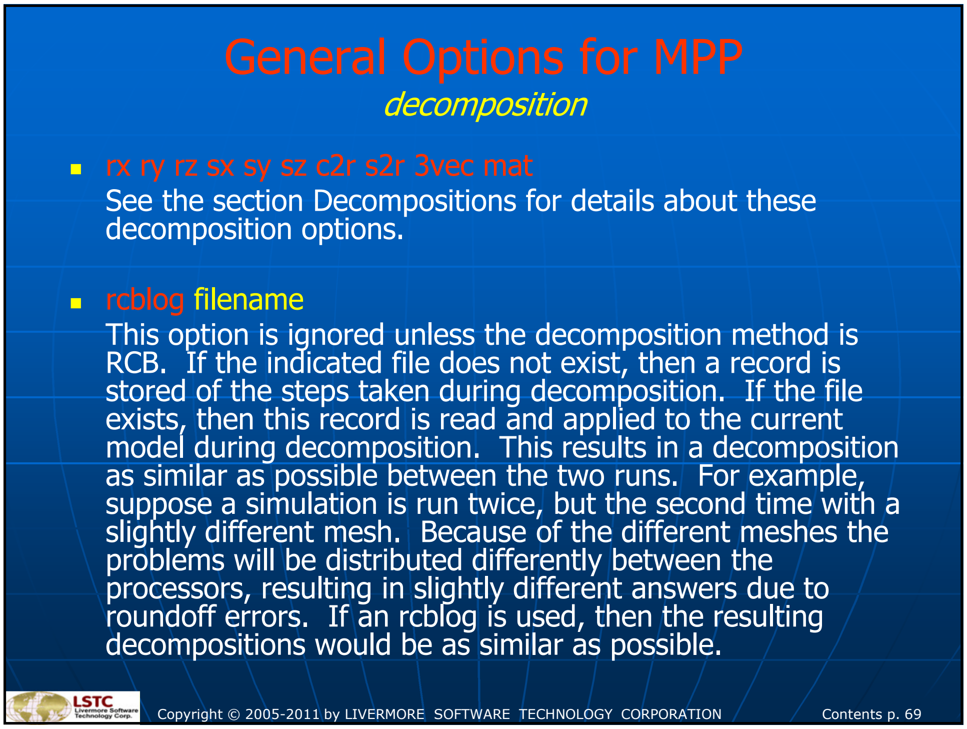
- local *path*

Path to a processor specific local directory for scratch/local files. This directory will be created if necessary. This is of primary use on systems where each processor has a local disk attached to it. Default = global path

- rep path

- transfer_files

Move output files back from local disk to starting directory or move restart files from starting directory to target local disk

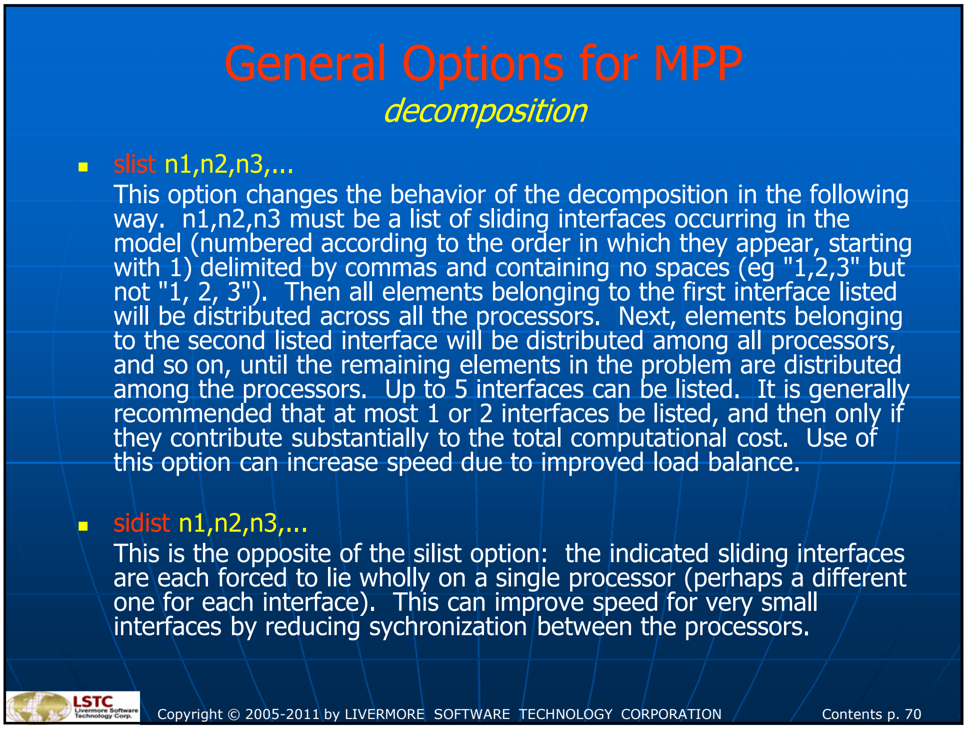


General Options for MPP

decomposition

- rx ry rz sx sy sz c2r s2r 3vec mat
  See the section Decompositions for details about these decomposition options.

- rcblog filename
  This option is ignored unless the decomposition method is RCB. If the indicated file does not exist, then a record is stored of the steps taken during decomposition. If the file exists, then this record is read and applied to the current model during decomposition. This results in a decomposition as similar as possible between the two runs. For example, suppose a simulation is run twice, but the second time with a slightly different mesh. Because of the different meshes the problems will be distributed differently between the processors, resulting in slightly different answers due to roundoff errors. If an rcblog is used, then the resulting decompositions would be as similar as possible.

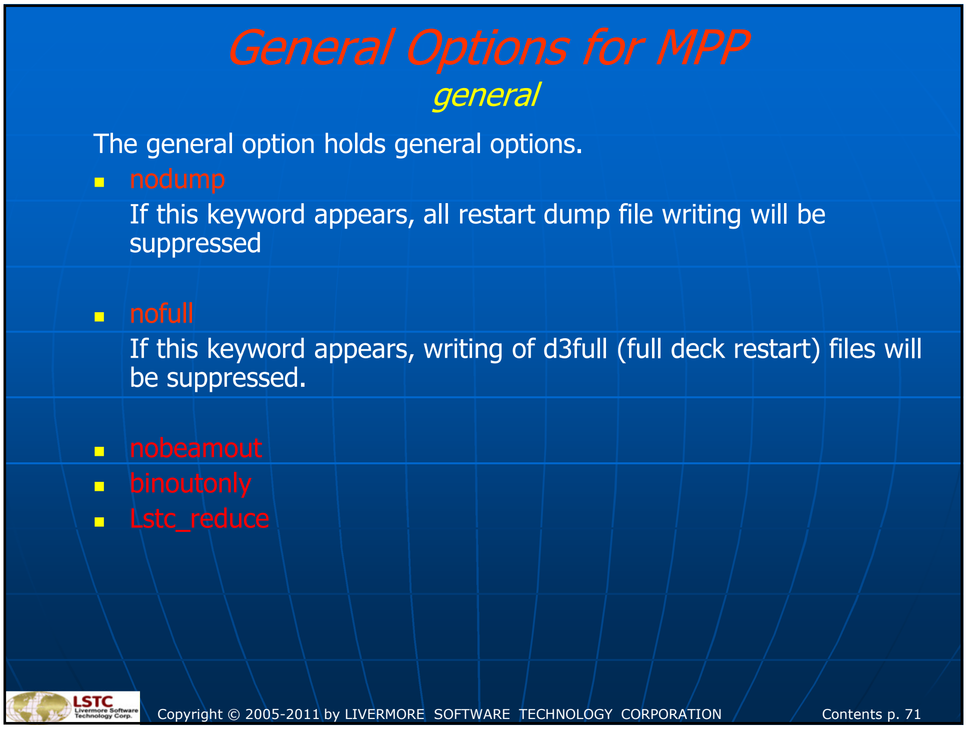


General Options for MPP

decomposition

- slist n1,n2,n3,...
  This option changes the behavior of the decomposition in the following way. n1,n2,n3 must be a list of sliding interfaces occurring in the model (numbered according to the order in which they appear, starting with 1) delimited by commas and containing no spaces (eg "1,2,3" but not "1, 2, 3"). Then all elements belonging to the first interface listed will be distributed across all the processors. Next, elements belonging to the second listed interface will be distributed among all processors, and so on, until the remaining elements in the problem are distributed among the processors. Up to 5 interfaces can be listed. It is generally recommended that at most 1 or 2 interfaces be listed, and then only if they contribute substantially to the total computational cost. Use of this option can increase speed due to improved load balance.

- sidist n1,n2,n3,...
  This is the opposite of the silist option: the indicated sliding interfaces are each forced to lie wholly on a single processor (perhaps a different one for each interface). This can improve speed for very small interfaces by reducing sychronization between the processors.



General Options for MPP

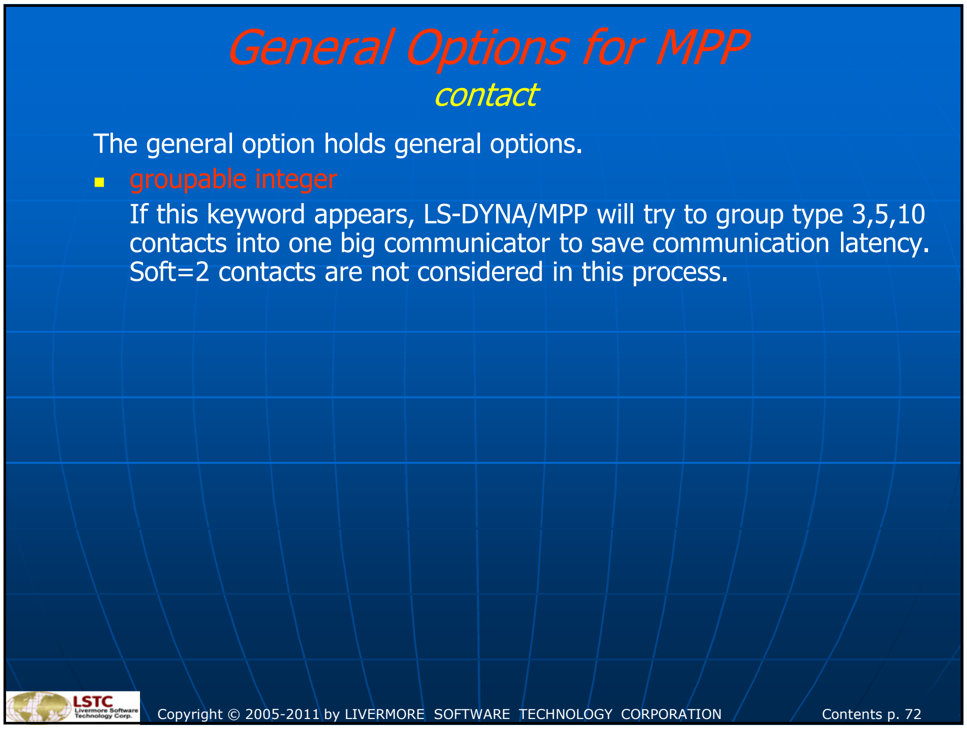
general

The general option holds general options.

- nodump
  If this keyword appears, all restart dump file writing will be suppressed

- nofull
  If this keyword appears, writing of d3full (full deck restart) files will be suppressed.

- nobeamout
- binoutonly
- Lstc_reduce



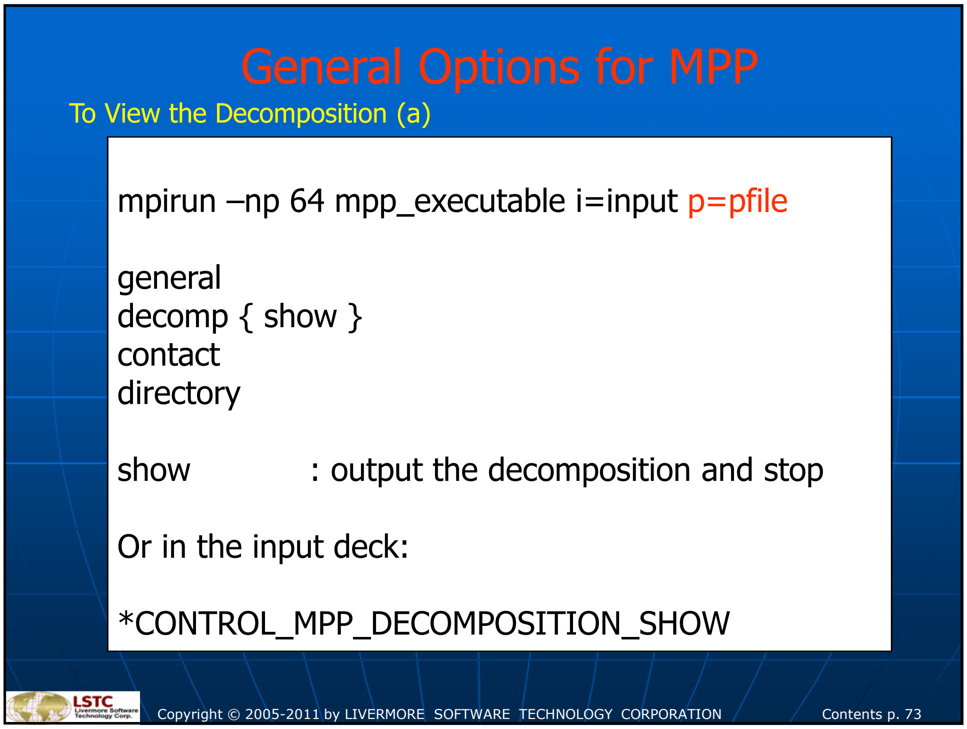
General Options for MPP

contact

The general option holds general options.

- groupable integer

  If this keyword appears, LS-DYNA/MPP will try to group type 3,5,10 contacts into one big communicator to save communication latency. Soft=2 contacts are not considered in this process.



General Options for MPP

To View the Decomposition (a)

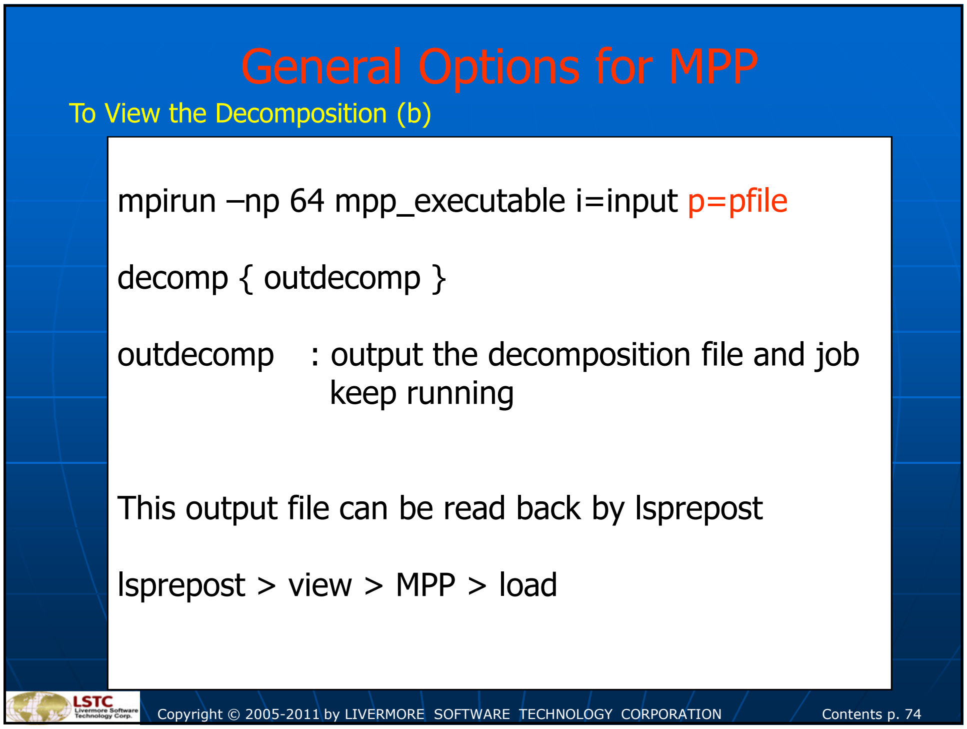
```
mpirun –np 64 mpp_executable i=input p=pfile

general
decomp { show }
contact
directory

show            : output the decomposition and stop

Or in the input deck:

*CONTROL_MPP_DECOMPOSITION_SHOW
```



General Options for MPP

To View the Decomposition (b)

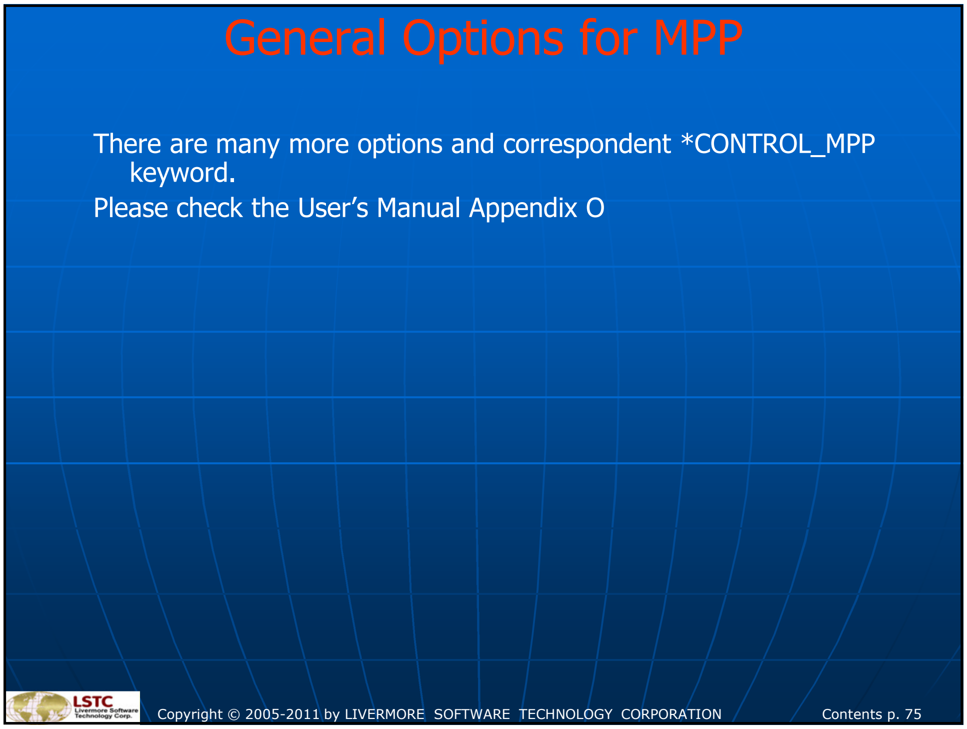
```
mpirun –np 64 mpp_executable i=input p=pfile
```

```
decomp { outdecomp }
```

outdecomp : output the decomposition file and job keep running

This output file can be read back by lsprepost

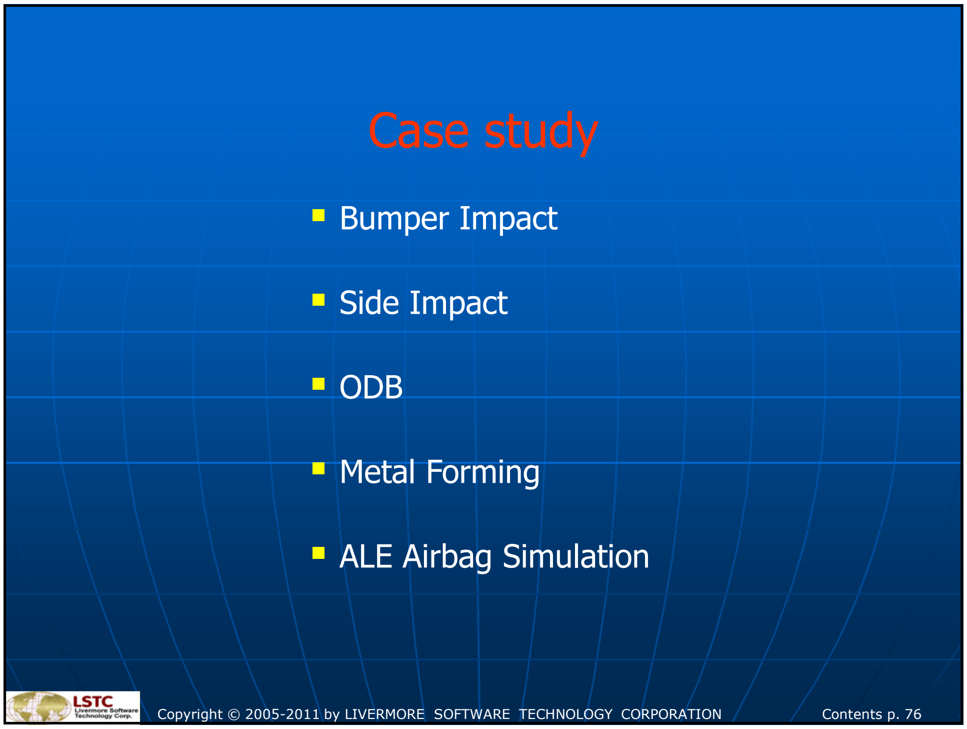
lsprepost > view > MPP > load



General Options for MPP

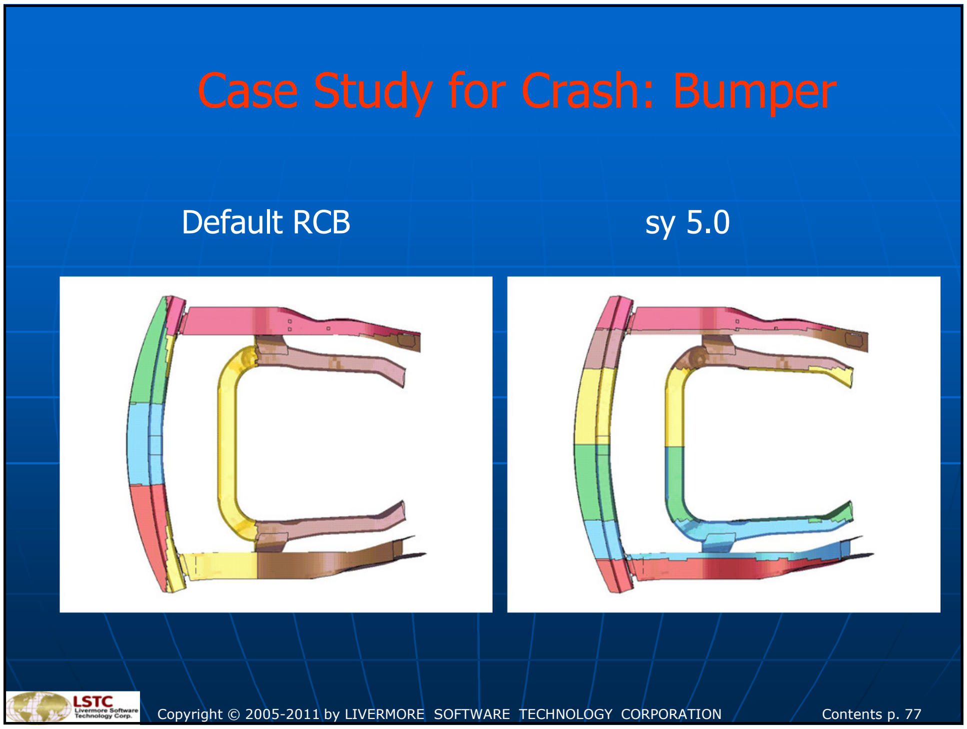
There are many more options and correspondent *CONTROL_MPP keyword.

Please check the User’s Manual Appendix O



Case study

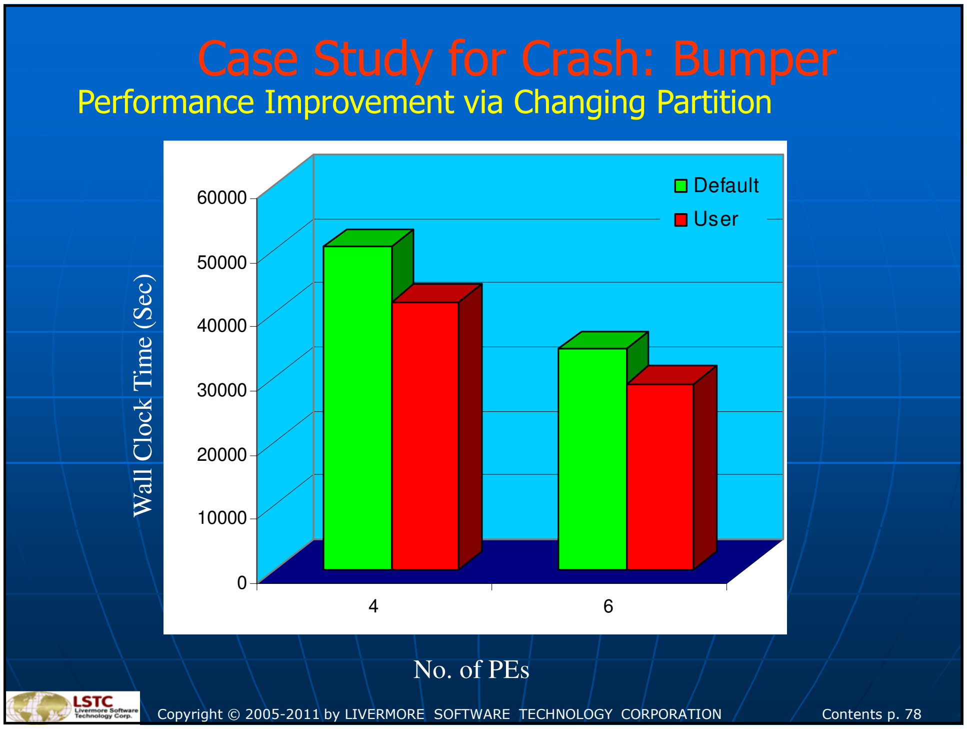
- Bumper Impact
- Side Impact
- ODB
- Metal Forming
- ALE Airbag Simulation



Case Study for Crash: Bumper

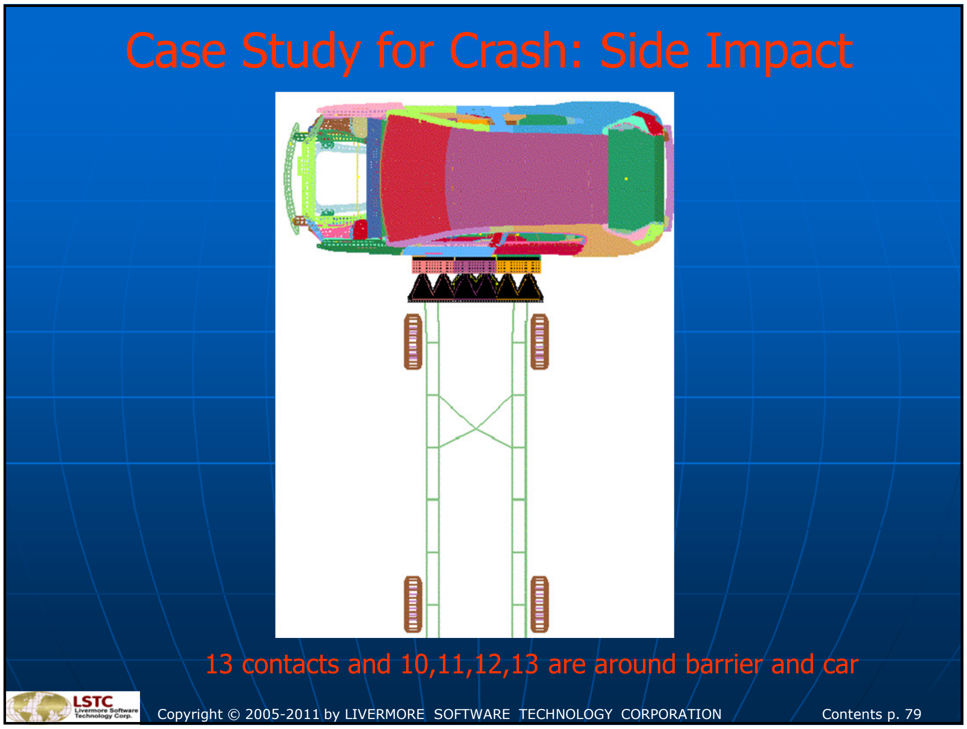
Default RCB

sy 5.0



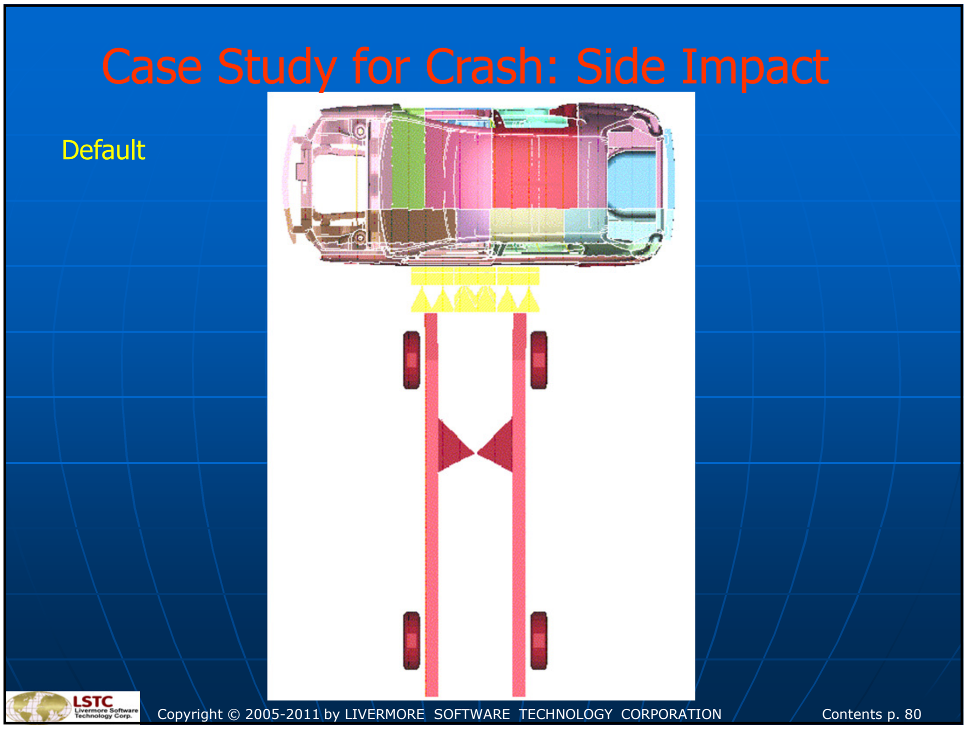
Case Study for Crash: Bumper

Performance Improvement via Changing Partition



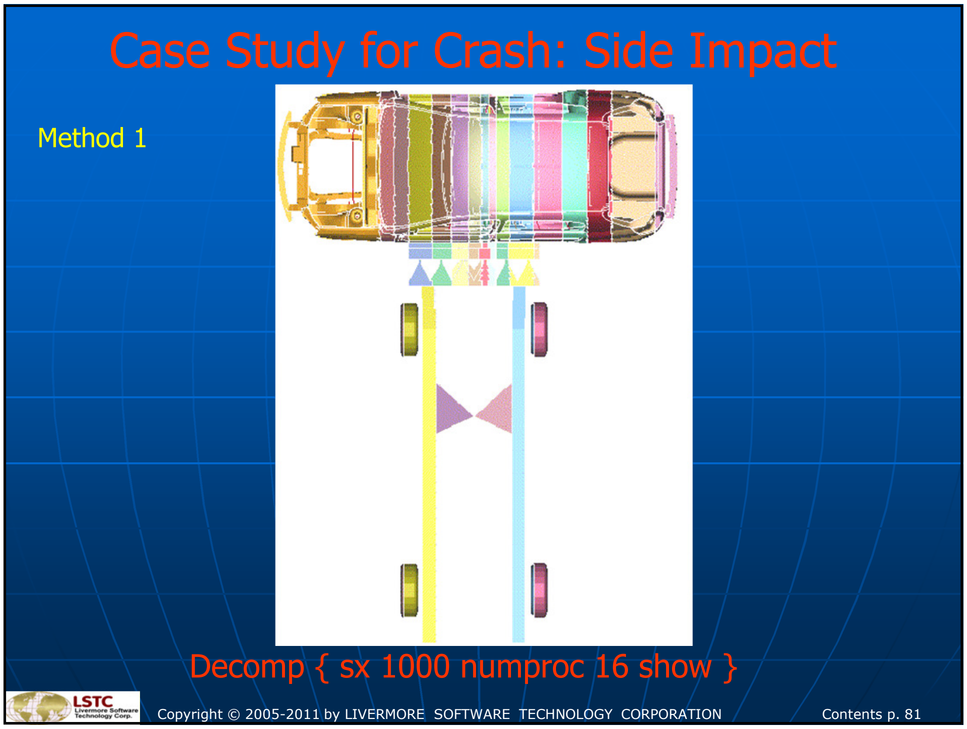
Case Study for Crash: Side Impact

13 contacts and 10,11,12,13 are around barrier and car



Case Study for Crash: Side Impact

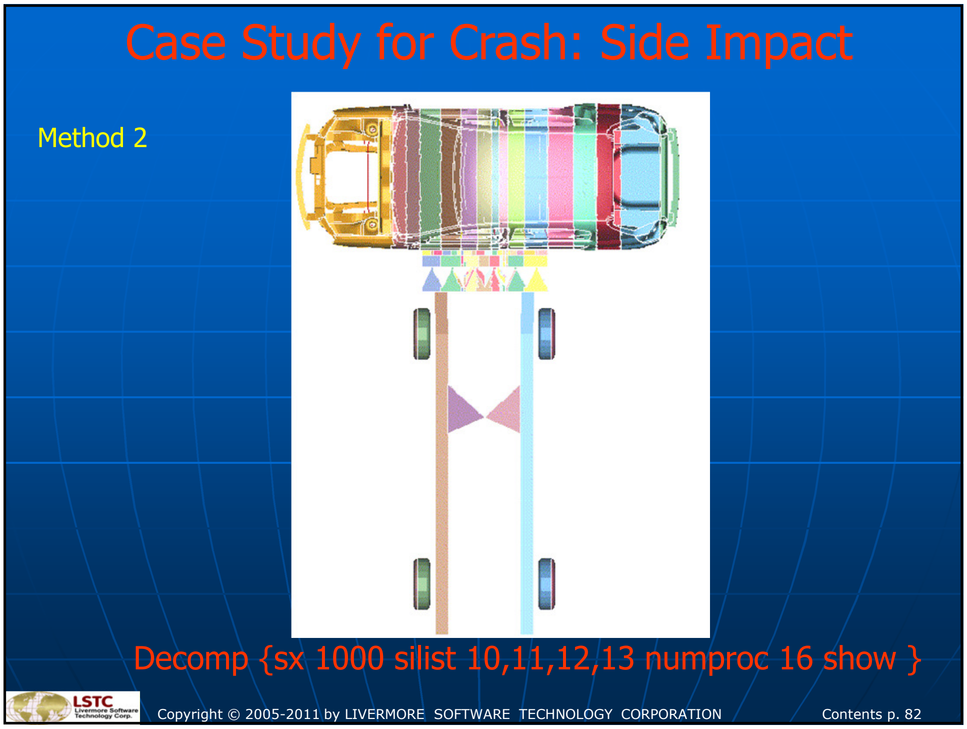
Default



Case Study for Crash: Side Impact

Method 1

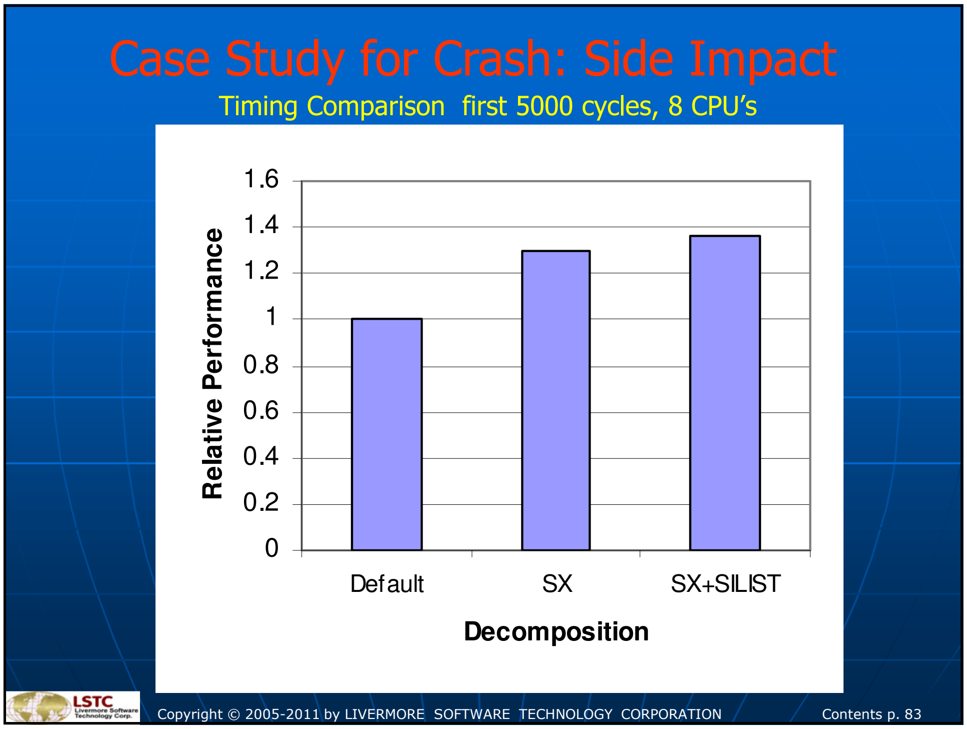
Decomp { sx 1000 numproc 16 show }



Case Study for Crash: Side Impact

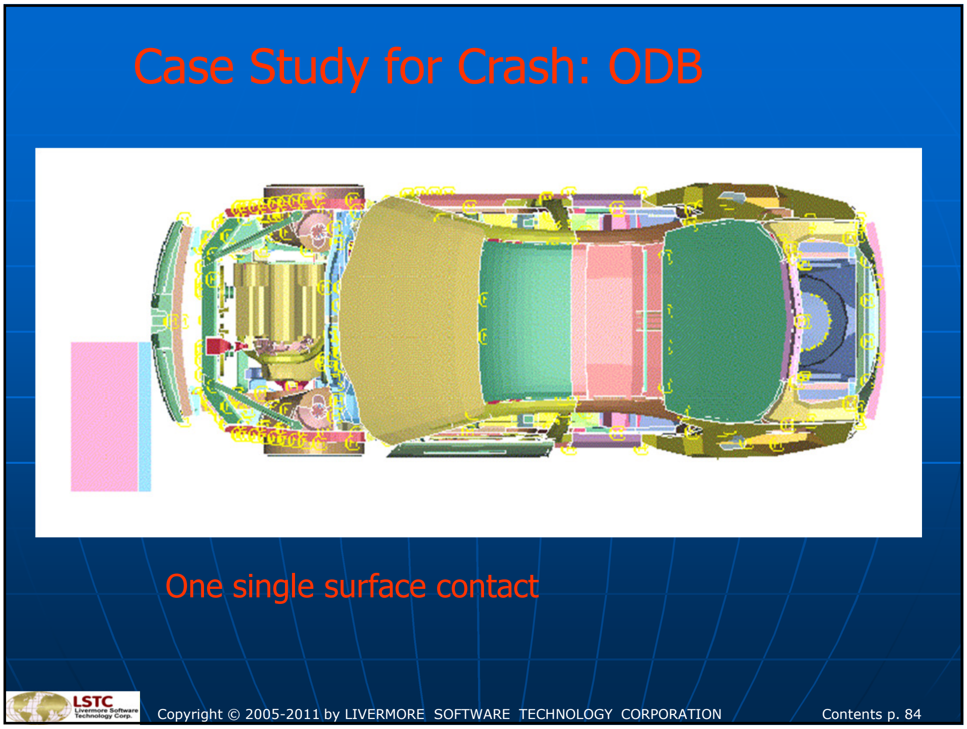
Method 2

Decomp {sx 1000 silist 10,11,12,13 numproc 16 show }



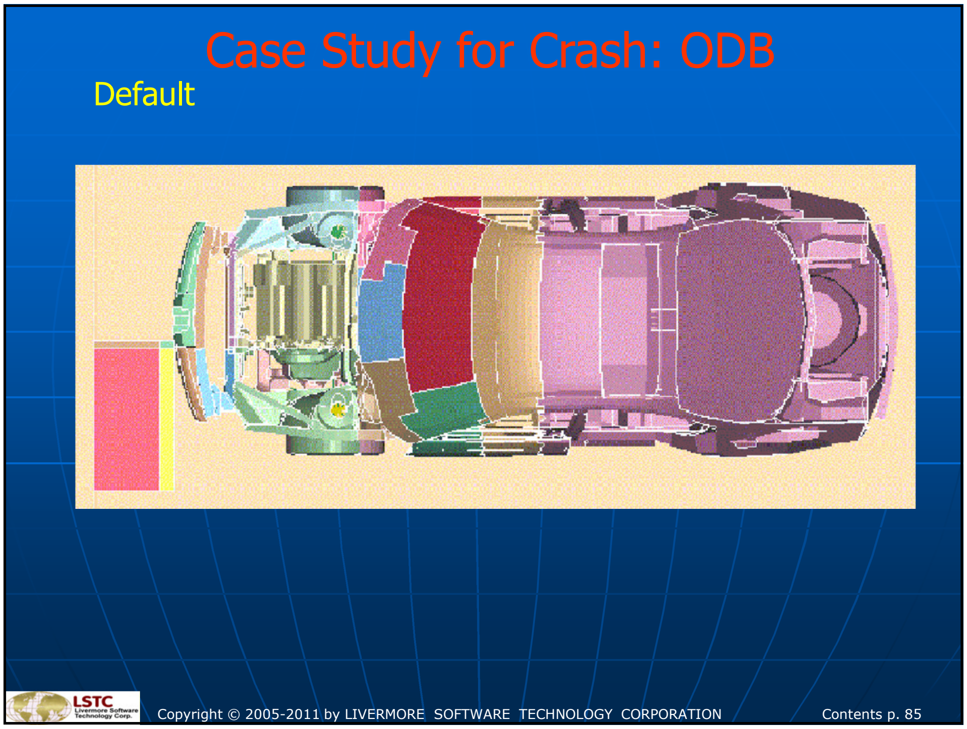
Case Study for Crash: Side Impact

Timing Comparison first 5000 cycles, 8 CPU's



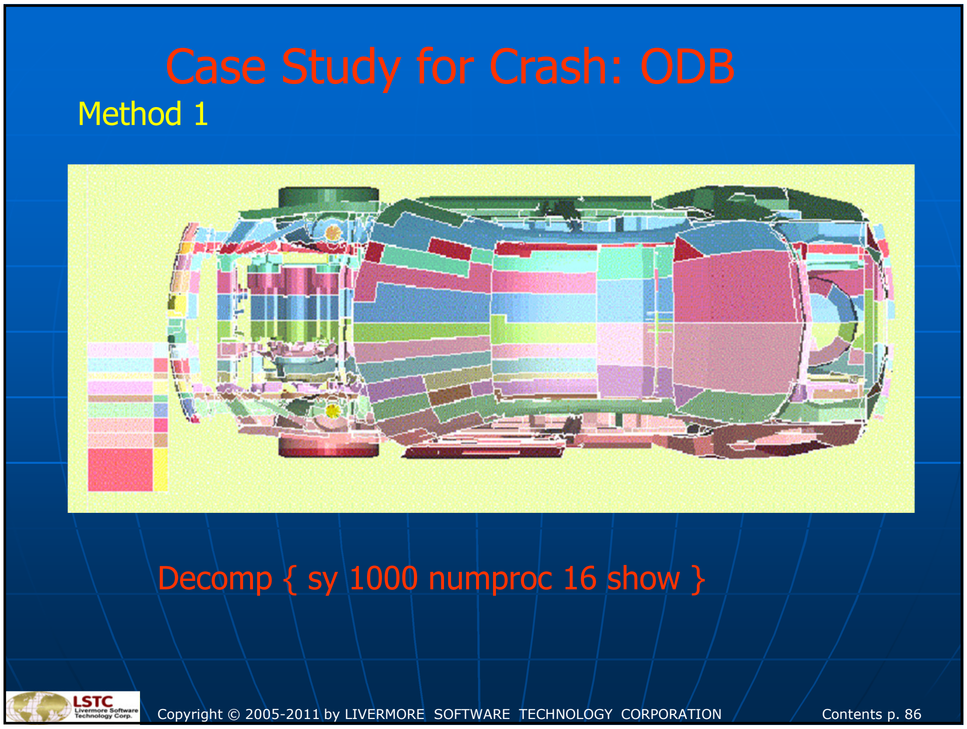
Case Study for Crash: ODB

One single surface contact



Case Study for Crash: ODB

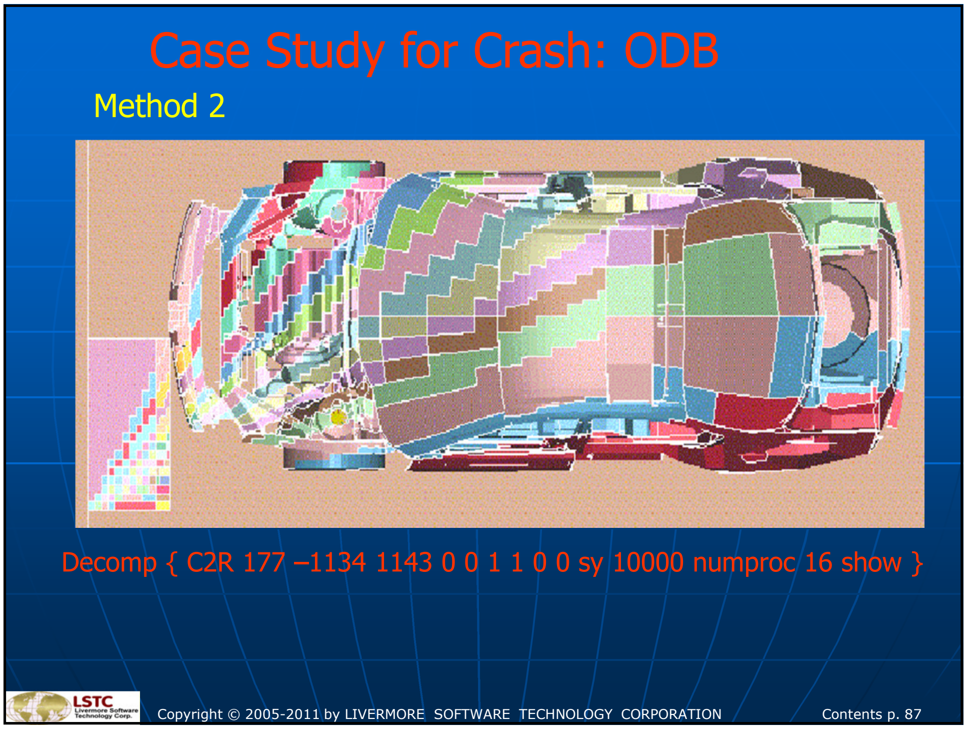
Default



Case Study for Crash: ODB

Method 1

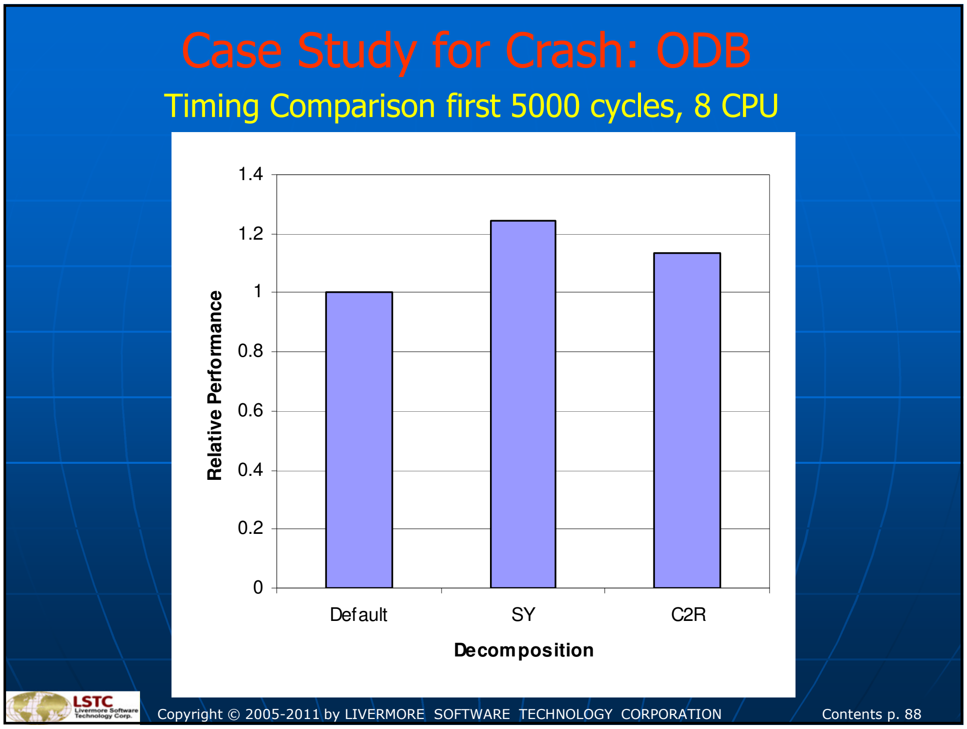
Decomp { sy 1000 numproc 16 show }



Case Study for Crash: ODB

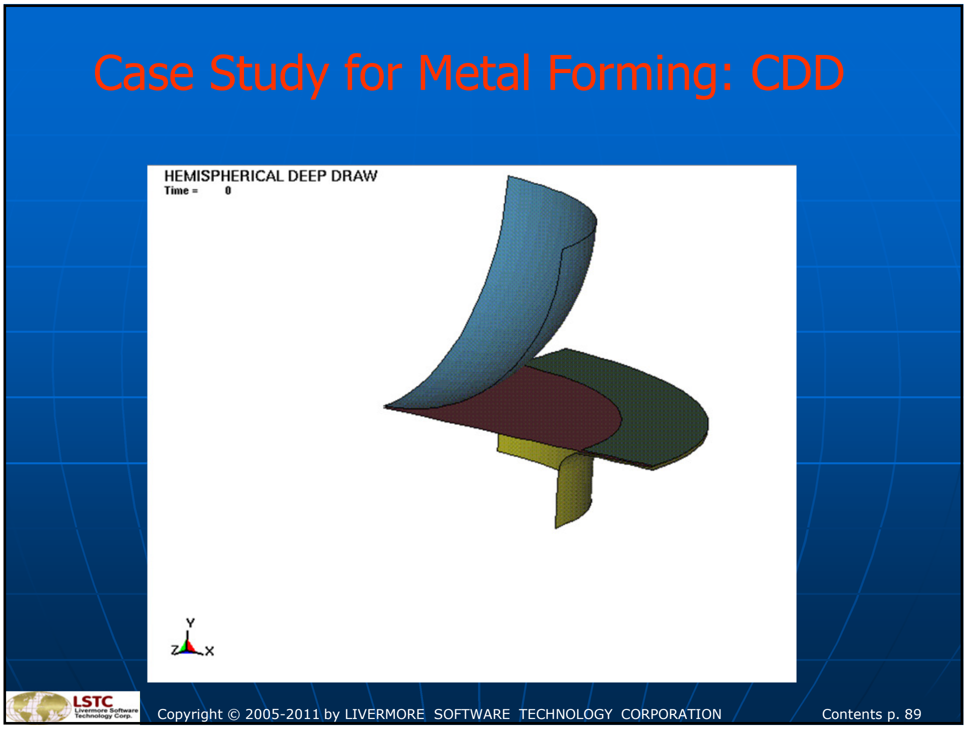
Method 2

Decomp { C2R 177 –1134 1143 0 0 1 1 0 0 sy 10000 numproc 16 show }

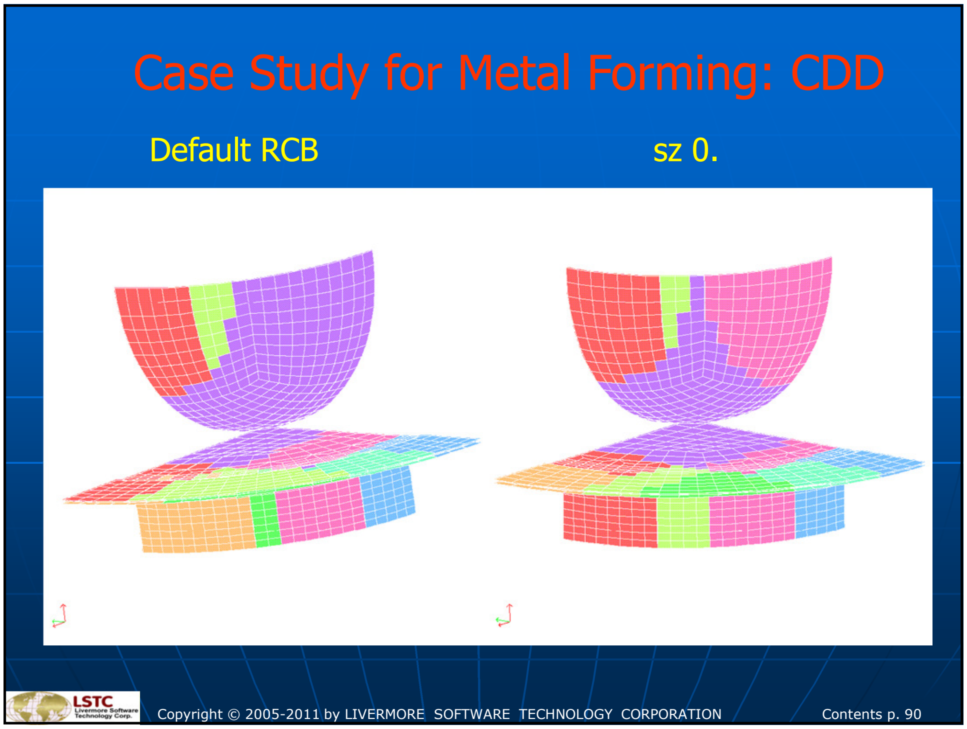


Case Study for Crash: ODB

Timing Comparison first 5000 cycles, 8 CPU



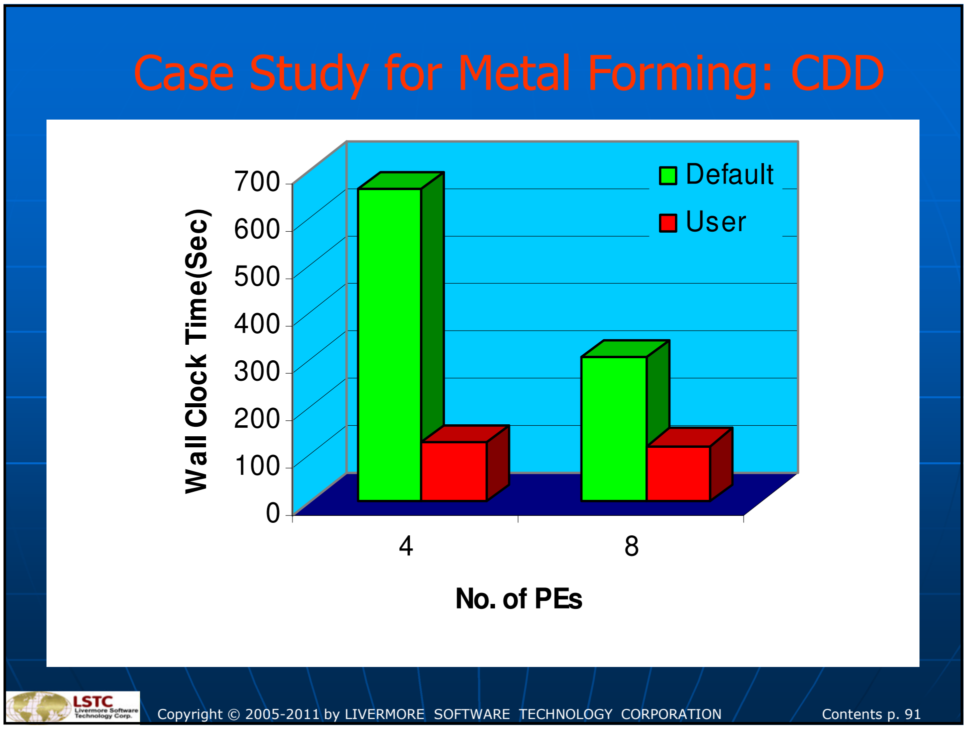
Case Study for Metal Forming: CDD



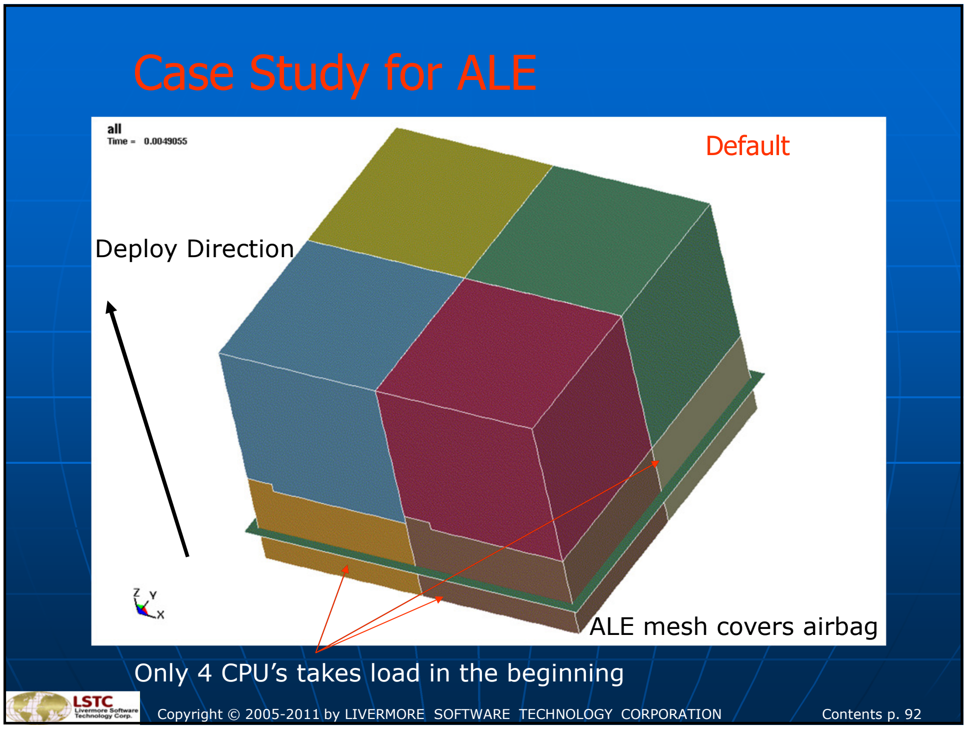
Case Study for Metal Forming: CDD

Default RCB

sz 0.



Case Study for Metal Forming: CDD



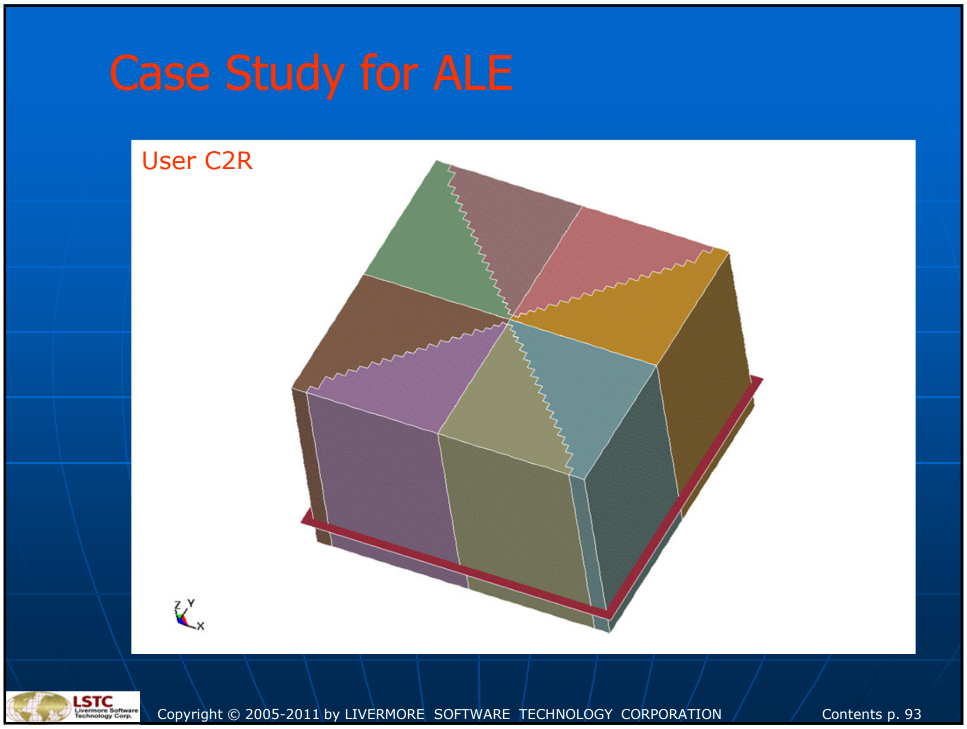
Case Study for ALE

Default

Deploy Direction

ALE mesh covers airbag

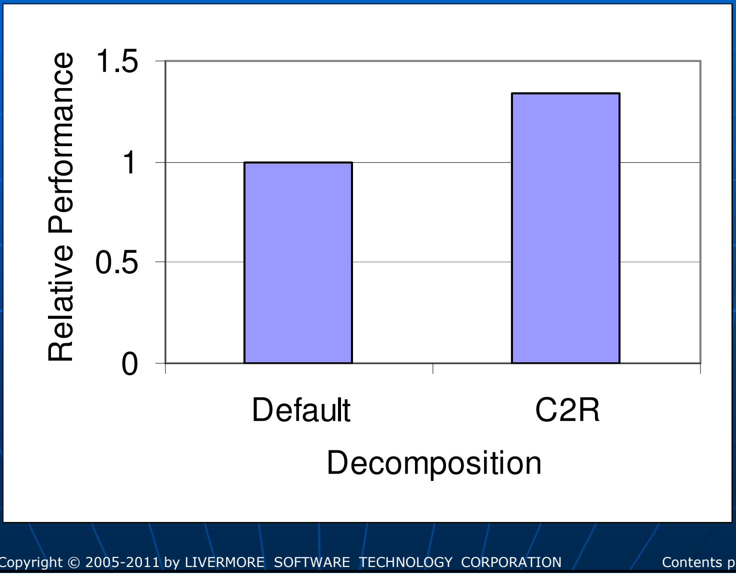
Only 4 CPU’s takes load in the beginning



Case Study for ALE

User C2R

ALE Airbag Timing Comparison
first 5000 cycles, 8 CPU



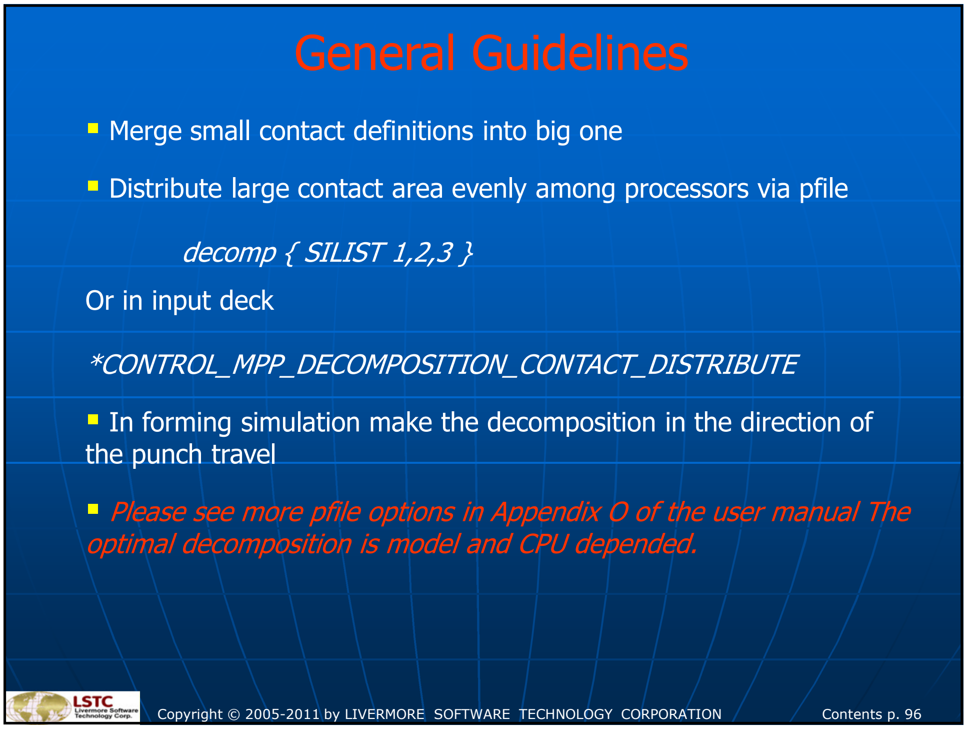
General Guidelines

- Use local file system "dir { local path }" if possible
  This allows MPP job to have scalable IO bandwidth

- Store end results via "dir { rep path }" to the share file system
  The files are moved through MPI calls which has higher bandwidth than NFS file system

- Distribute expansive features or elements to all processors
  i.e. CPM airbag, ALE elements, SPH elements, etc
  (*CONTROL_MPP_DECOMPOSITION_BAGREF)
  (*CONTROL_MPP_DECOMPOSITION_ARRANGE_PARTS)
  (*CONTROL_MPP_DECOMPOSITION_DISTRIBUTE_ALE_ELE, etc)

- For number of processors < 16, try to partition model along the direction of initial velocity (use e.g. automatic decomposition (*CONTROL_MPP_DECOMPOSITION_AUTO)



General Guidelines

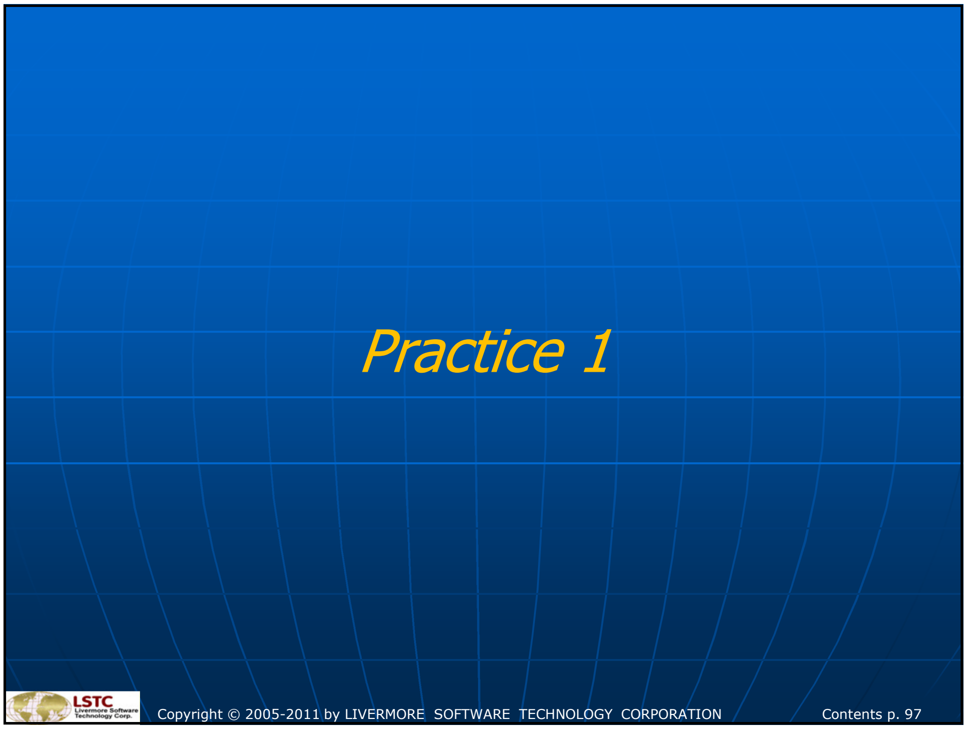
- Merge small contact definitions into big one
- Distribute large contact area evenly among processors via pfile

decomp { SILIST 1,2,3 }

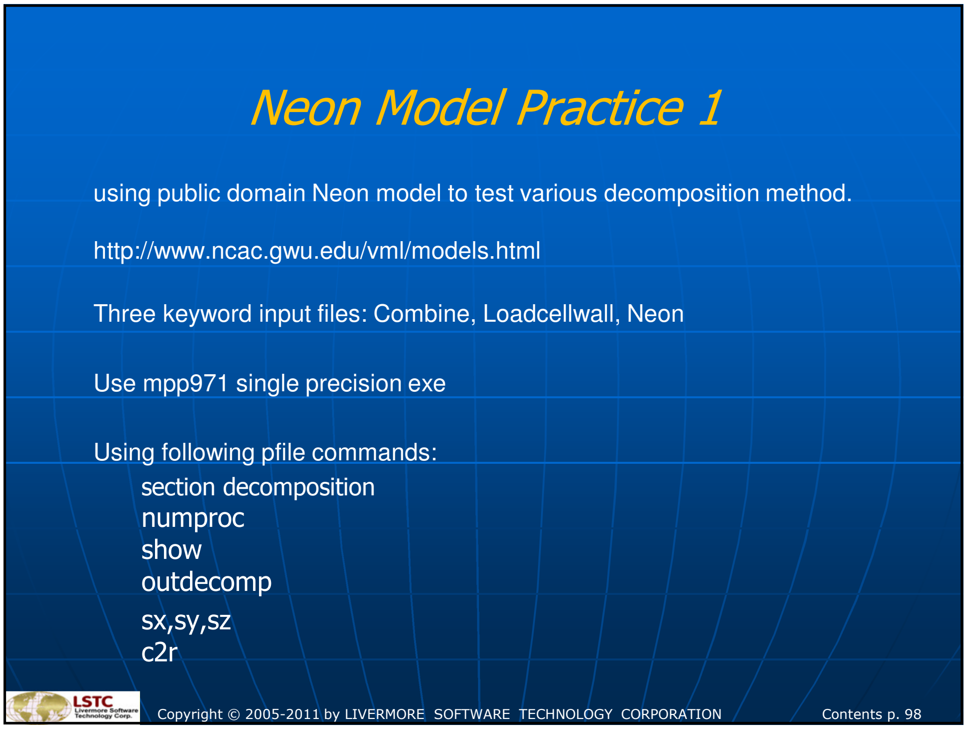
Or in input deck

**CONTROL_MPP_DECOMPOSITION_CONTACT_DISTRIBUTE*

- In forming simulation make the decomposition in the direction of the punch travel
- *Please see more pfile options in Appendix O of the user manual The optimal decomposition is model and CPU depended.*



Practice 1



Neon Model Practice 1

using public domain Neon model to test various decomposition method.

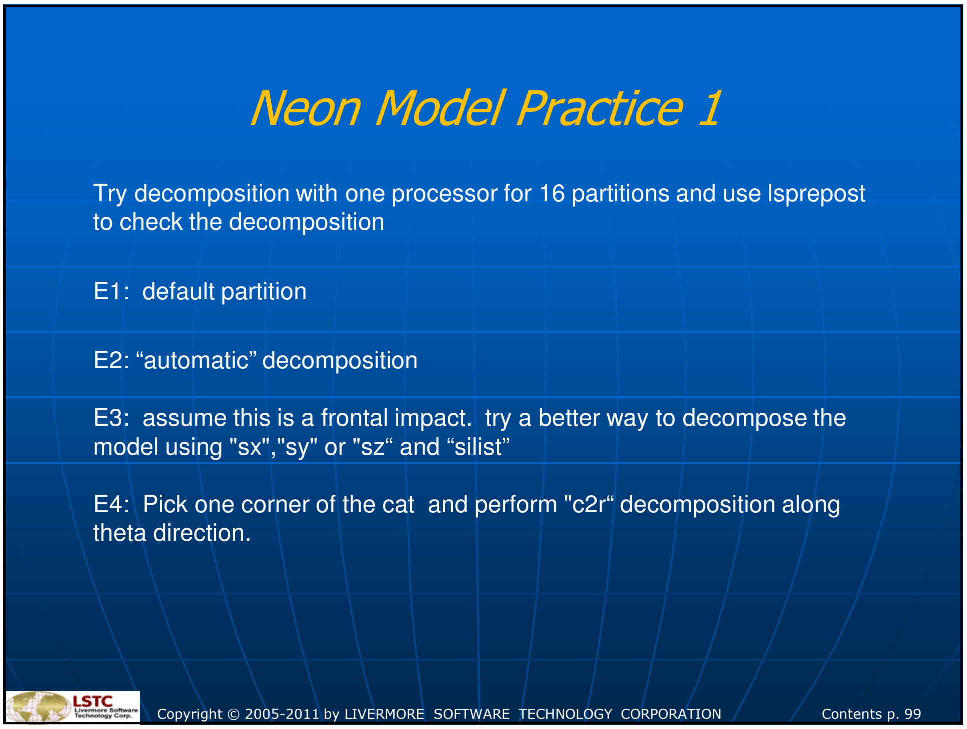
http://www.ncac.gwu.edu/vml/models.html

Three keyword input files: Combine, Loadcellwall, Neon

Use mpp971 single precision exe

Using following pfile commands:

section decomposition
numproc
show
outdecomp
sx,sy,sz
c2r



Neon Model Practice 1

Try decomposition with one processor for 16 partitions and use lsprepost to check the decomposition

E1: default partition

E2: “automatic” decomposition

E3: assume this is a frontal impact. try a better way to decompose the model using "sx","sy" or "sz“ and “silist”

E4: Pick one corner of the cat and perform "c2r“ decomposition along theta direction.

Restart and Pre-decomposition

Restart

- Restart is in MPP-DYNA is different from LS-DYNA, The files are called d3dump##.xxxx or d3fulll##, where ## is a number.

Simple restart: mpirun –np 5 mpp970 r=d3dump09

MPP-DYNA finds the child files

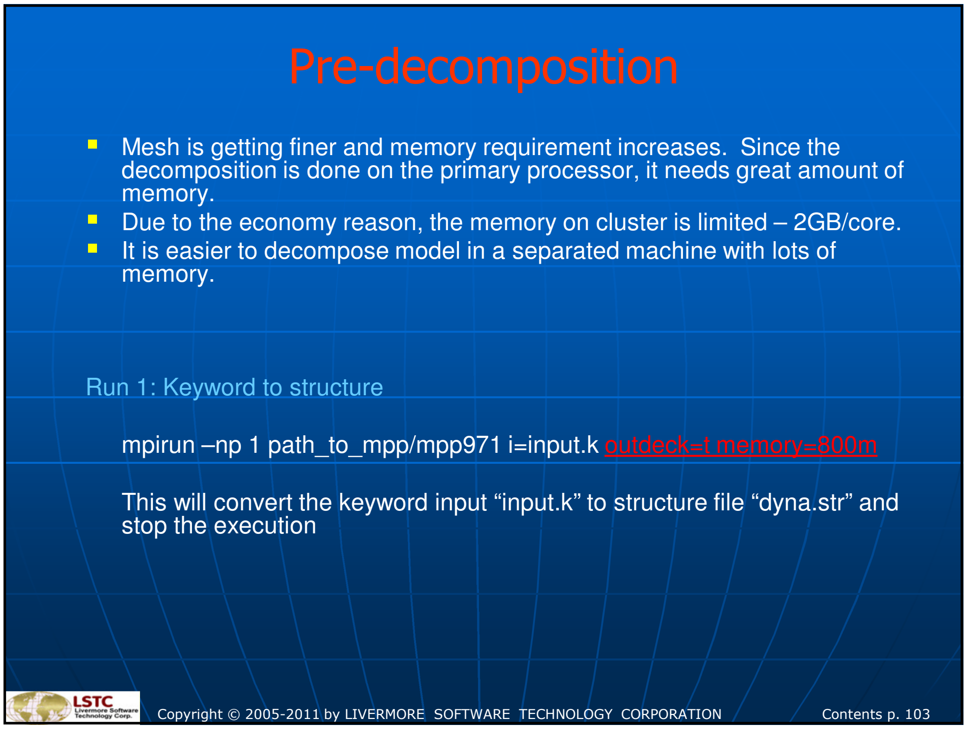
Small restart: mpirun –np 5 mpp970 i=small.k r=d3dump09

The small restart may have problems. If it does, please report it to LSTC and we will fix it.

Full restart: mpirun –np 5 mpp970 i=full.k n=d3full09

Remember *stress_initialization in the inputdeck
Can change ncpu in full restart
The full restart can have problems

- Since the Small and Full restart can give problems – check carefully the results



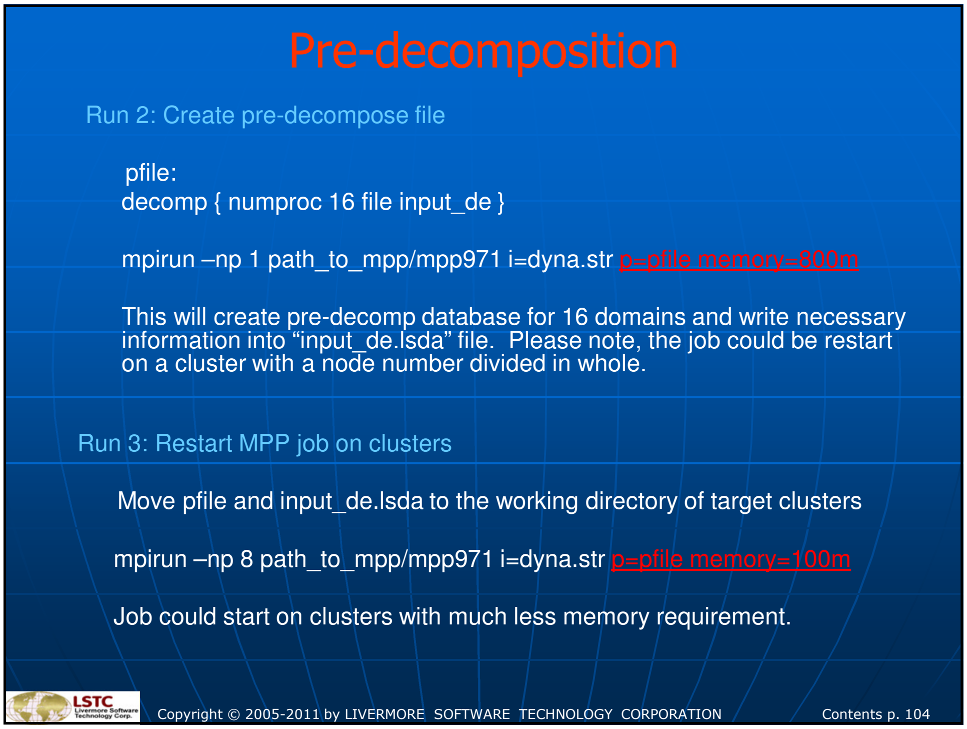
Pre-decomposition

- Mesh is getting finer and memory requirement increases. Since the decomposition is done on the primary processor, it needs great amount of memory.
- Due to the economy reason, the memory on cluster is limited – 2GB/core.
- It is easier to decompose model in a separated machine with lots of memory.

Run 1: Keyword to structure

```
mpirun –np 1 path_to_mpp/mpp971 i=input.k outdeck=t memory=800m
```

This will convert the keyword input “input.k” to structure file “dyna.str” and stop the execution



Pre-decomposition

Run 2: Create pre-decompose file

```
pfile:
decomp { numproc 16 file input_de }

mpirun –np 1 path_to_mpp/mpp971 i=dyna.str p=pfile memory=800m
```

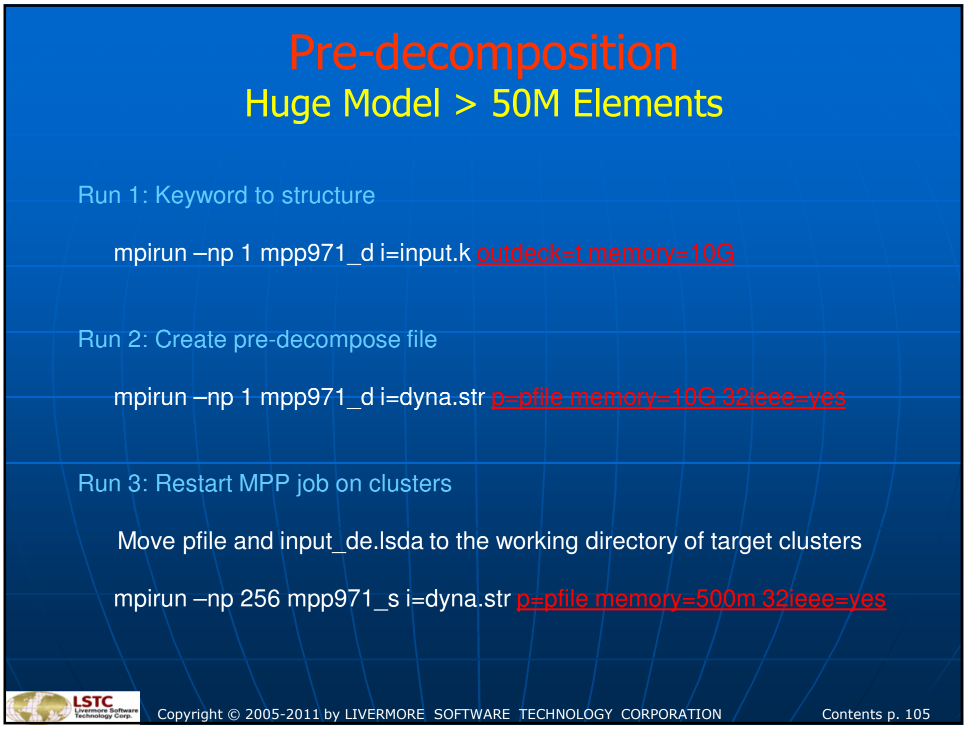
This will create pre-decomp database for 16 domains and write necessary information into “input_de.lsda” file. Please note, the job could be restart on a cluster with a node number divided in whole.

Run 3: Restart MPP job on clusters

Move pfile and input_de.lsda to the working directory of target clusters

```
mpirun –np 8 path_to_mpp/mpp971 i=dyna.str p=pfile memory=100m
```

Job could start on clusters with much less memory requirement.



Pre-decomposition

Huge Model > 50M Elements

Run 1: Keyword to structure

```
mpirun –np 1 mpp971_d i=input.k outdeck=t memory=10G
```

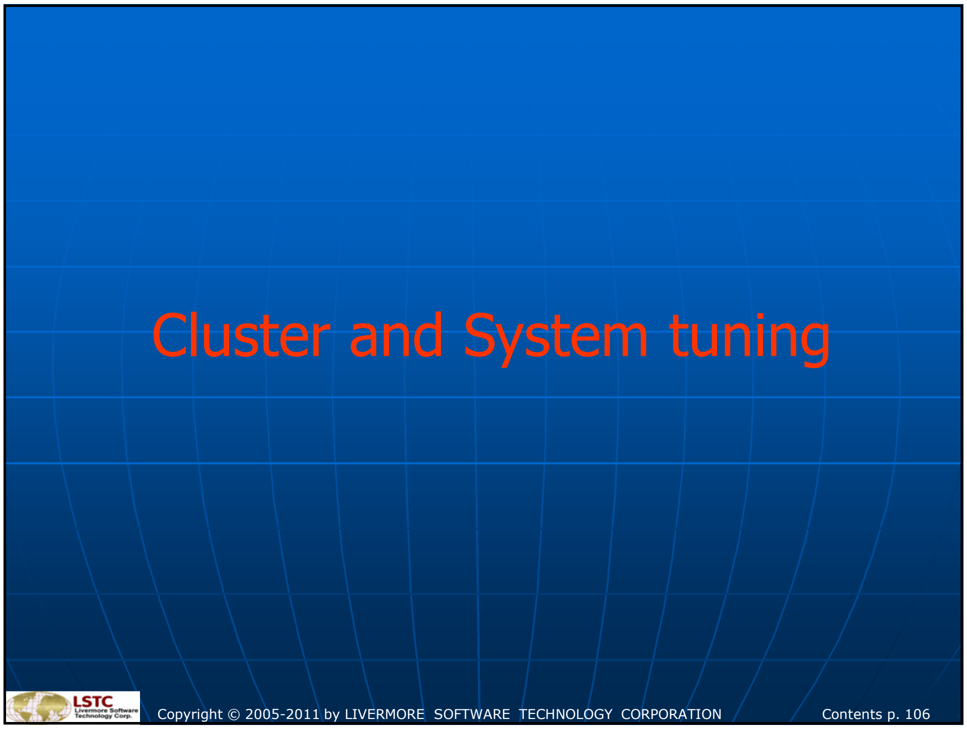
Run 2: Create pre-decompose file

```
mpirun –np 1 mpp971_d i=dyna.str p=pfile memory=10G 32ieee=yes
```

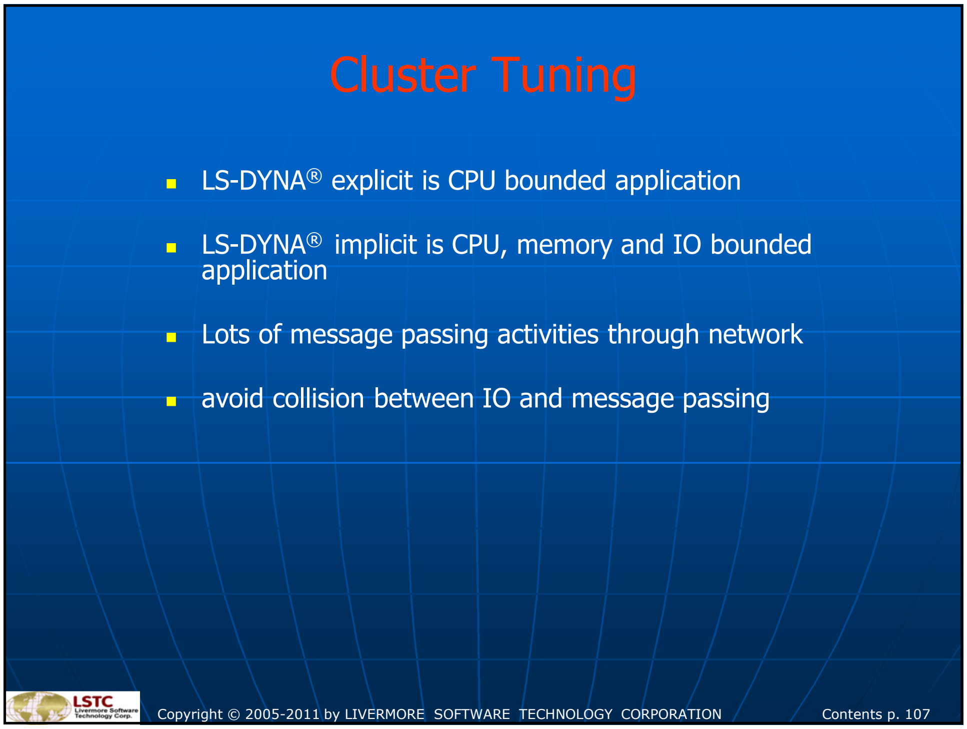
Run 3: Restart MPP job on clusters

Move pfile and input_de.lsda to the working directory of target clusters

```
mpirun –np 256 mpp971_s i=dyna.str p=pfile memory=500m 32ieee=yes
```

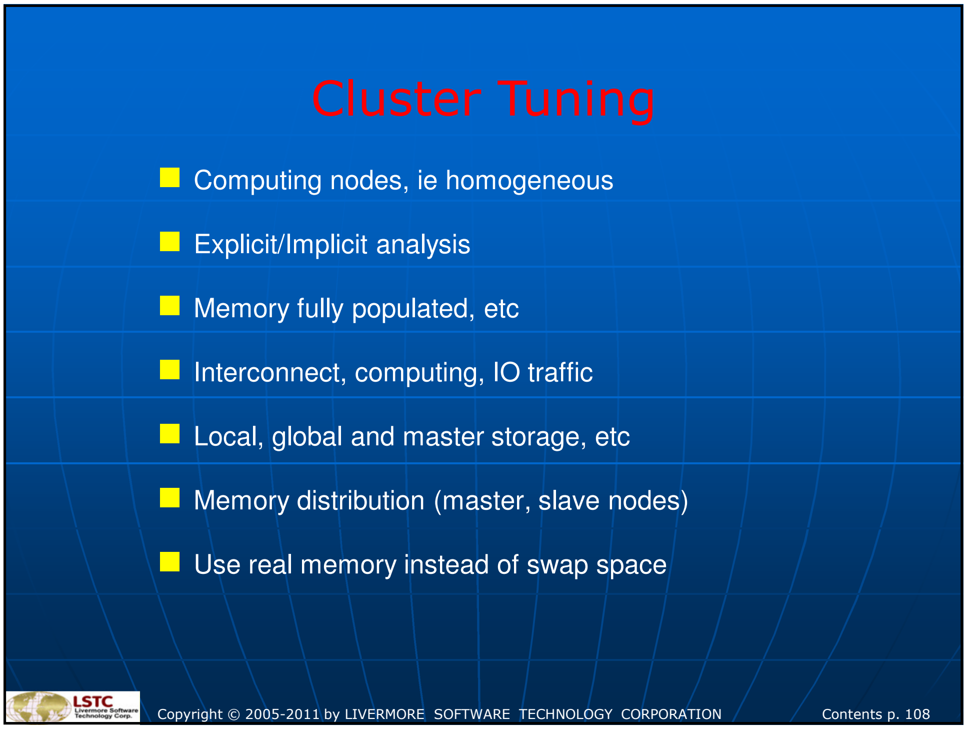


Cluster and System tuning



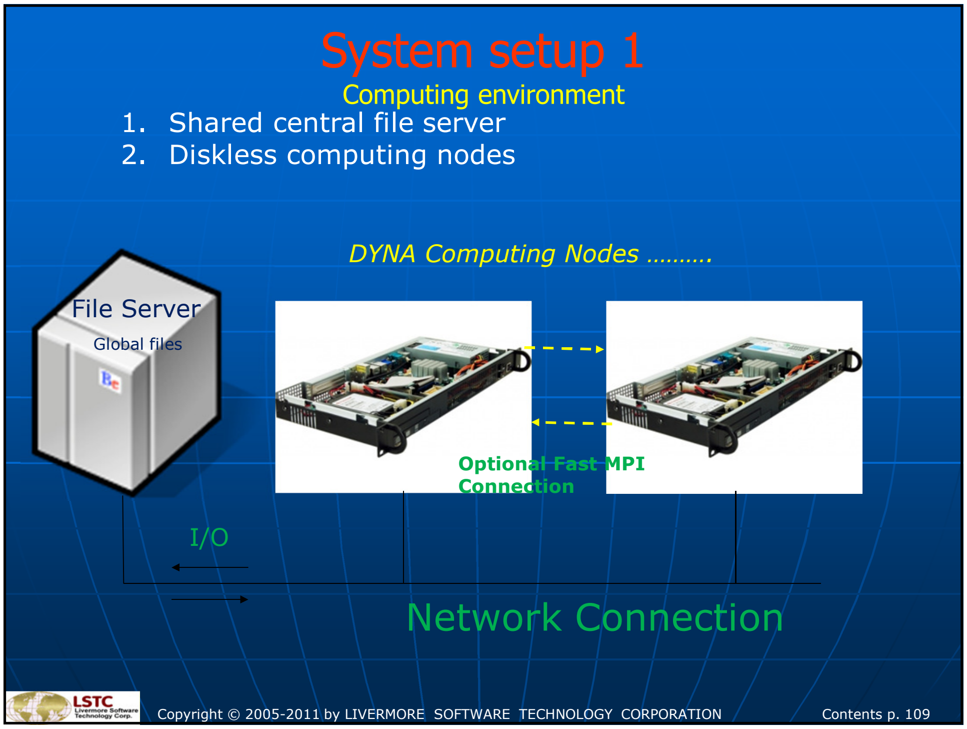
Cluster Tuning

- LS-DYNA® explicit is CPU bounded application
- LS-DYNA® implicit is CPU, memory and IO bounded application
- Lots of message passing activities through network
- avoid collision between IO and message passing



Cluster Tuning

- Computing nodes, ie homogeneous
- Explicit/Implicit analysis
- Memory fully populated, etc
- Interconnect, computing, IO traffic
- Local, global and master storage, etc
- Memory distribution (master, slave nodes)
- Use real memory instead of swap space



System setup 1

Computing environment

1. Shared central file server
2. Diskless computing nodes

DYNA Computing Nodes ……….

File Server

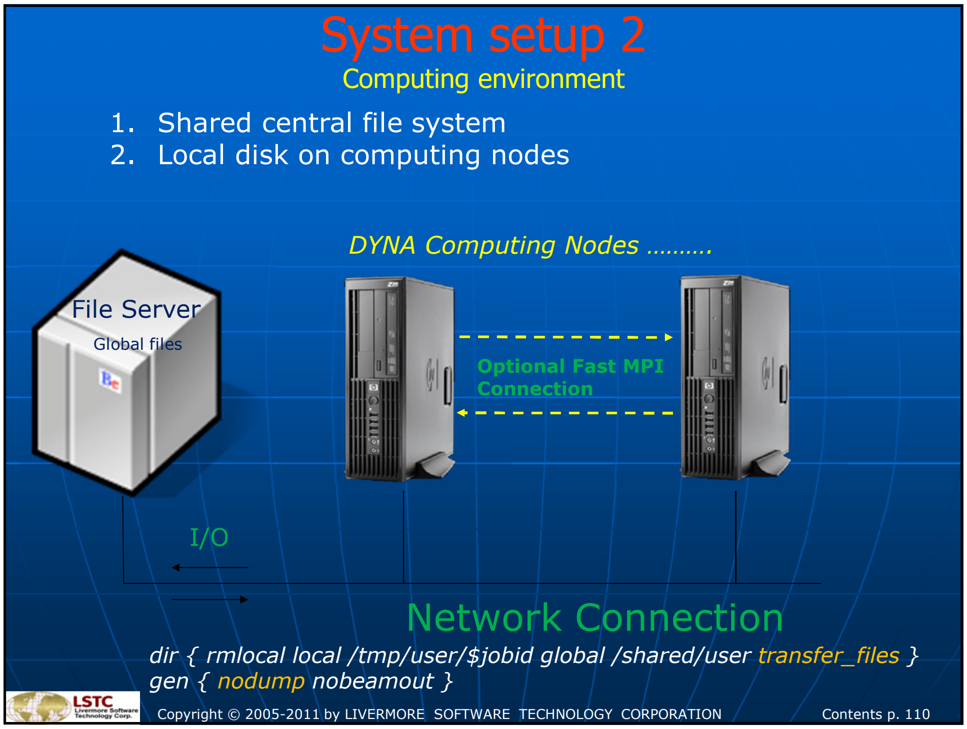
Global files

Optional Fast MPI Connection

I/O

Network Connection

Contents p. 109



System setup 2

Computing environment

1. Shared central file system
2. Local disk on computing nodes

DYNA Computing Nodes ……….

File Server

Global files

Optional Fast MPI Connection

I/O

Network Connection

dir { rmlocal local /tmp/user/$jobid global /shared/user transfer_files }
gen { nodump nobeamout }

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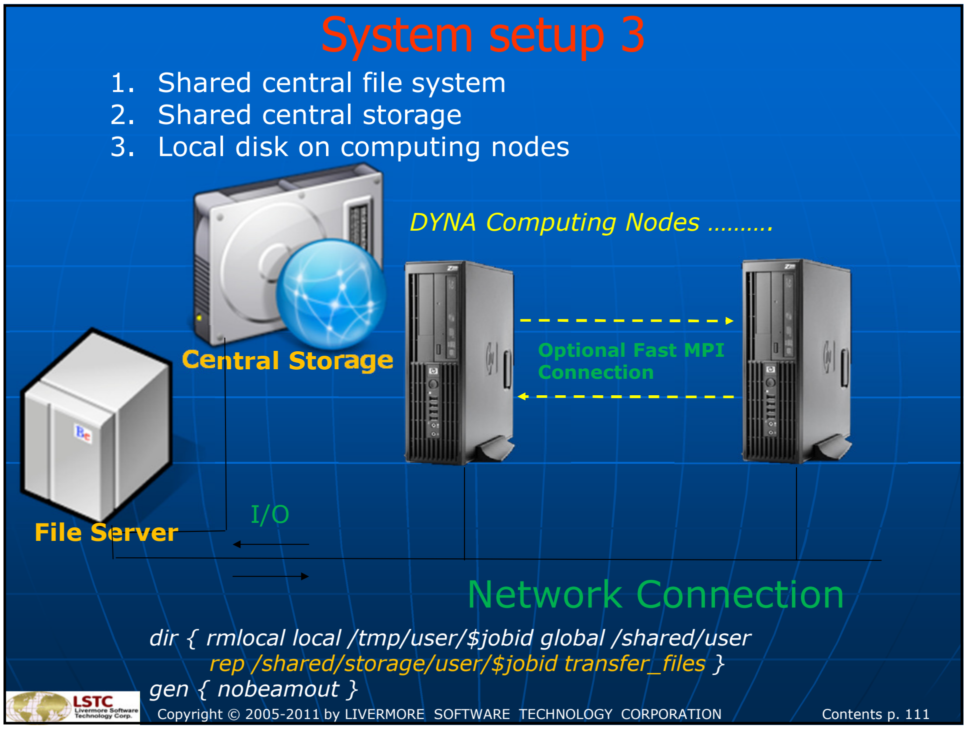


System setup 3

1. Shared central file system
2. Shared central storage
3. Local disk on computing nodes

DYNA Computing Nodes ……….

Central Storage

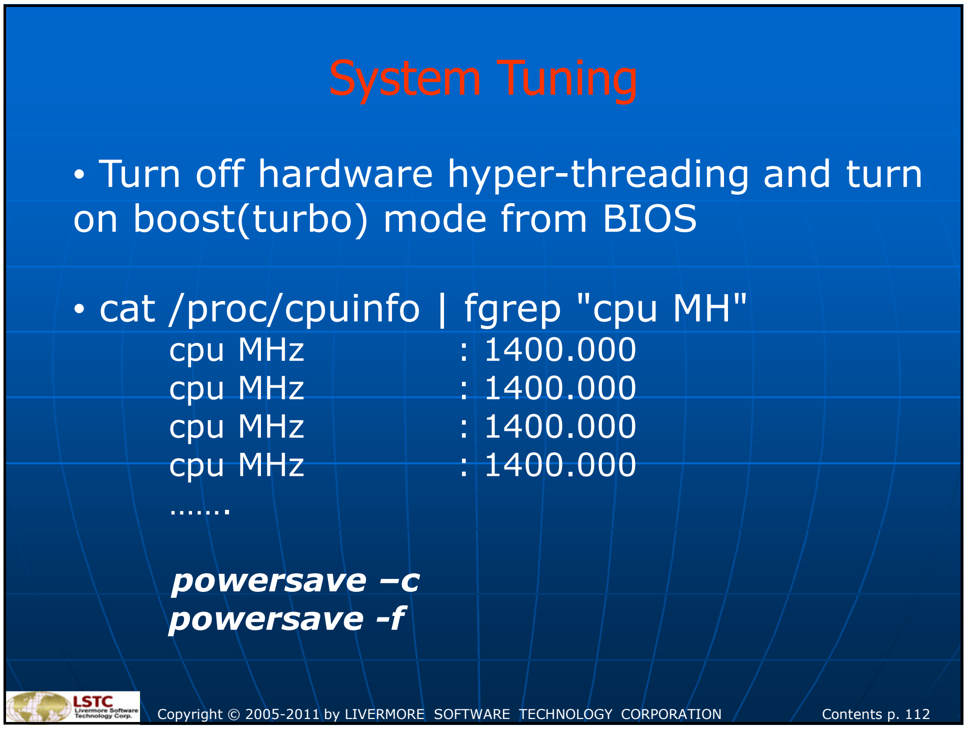
Optional Fast MPI Connection

File Server

I/O

Network Connection

dir { rmlocal local /tmp/user/$jobid global /shared/user
rep /shared/storage/user/$jobid transfer_files }
gen { nobeamout }



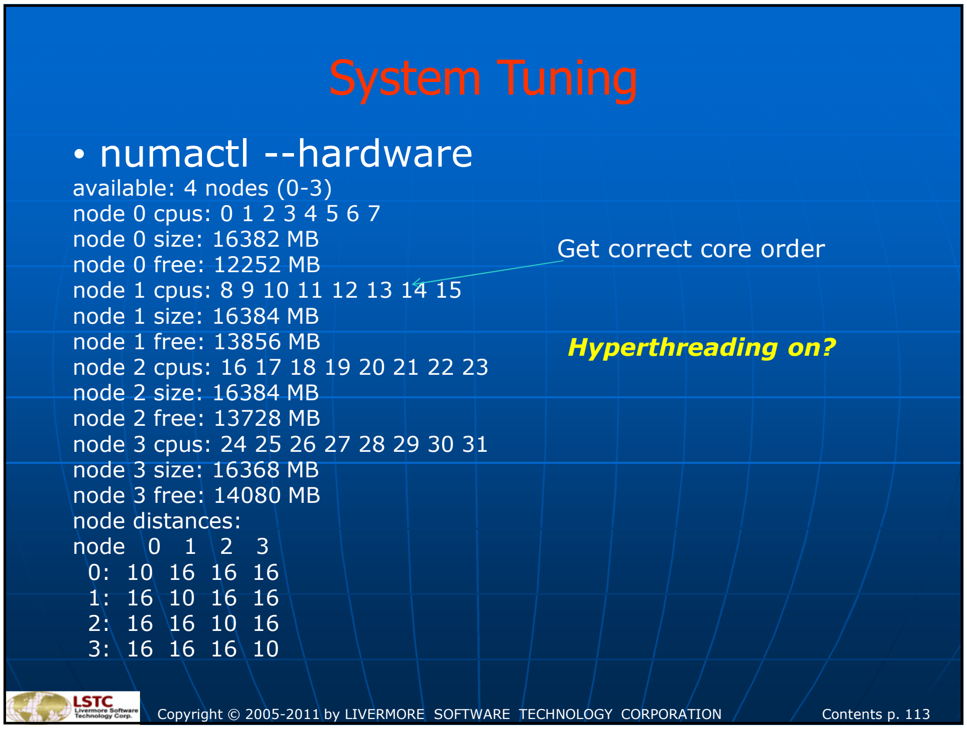
System Tuning

- Turn off hardware hyper-threading and turn on boost(turbo) mode from BIOS

- cat /proc/cpuinfo | fgrep "cpu MH"

```
cpu MHz          : 1400.000
cpu MHz          : 1400.000
cpu MHz          : 1400.000
cpu MHz          : 1400.000
…….
```

powersave –c
powersave -f



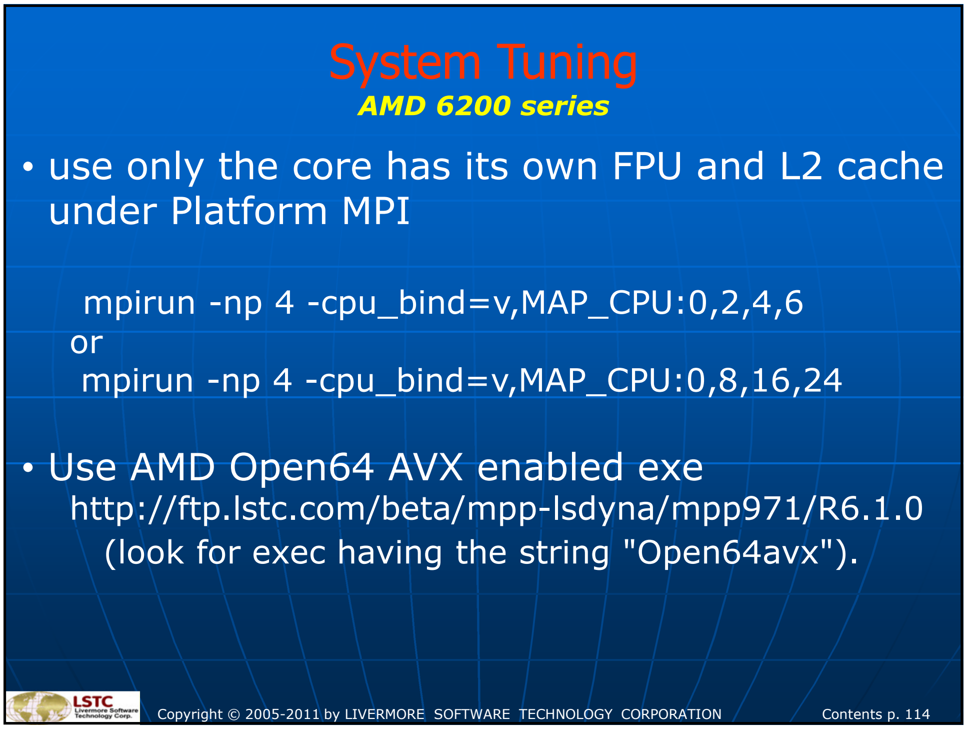
System Tuning

- numactl --hardware

```
available: 4 nodes (0-3)
node 0 cpus: 0 1 2 3 4 5 6 7
node 0 size: 16382 MB
node 0 free: 12252 MB
node 1 cpus: 8 9 10 11 12 13 14 15
node 1 size: 16384 MB
node 1 free: 13856 MB
node 2 cpus: 16 17 18 19 20 21 22 23
node 2 size: 16384 MB
node 2 free: 13728 MB
node 3 cpus: 24 25 26 27 28 29 30 31
node 3 size: 16368 MB
node 3 free: 14080 MB
node distances:
node   0   1   2   3
  0:  10  16  16  16
  1:  16  10  16  16
  2:  16  16  10  16
  3:  16  16  16  10
```

Get correct core order

Hyperthreading on?



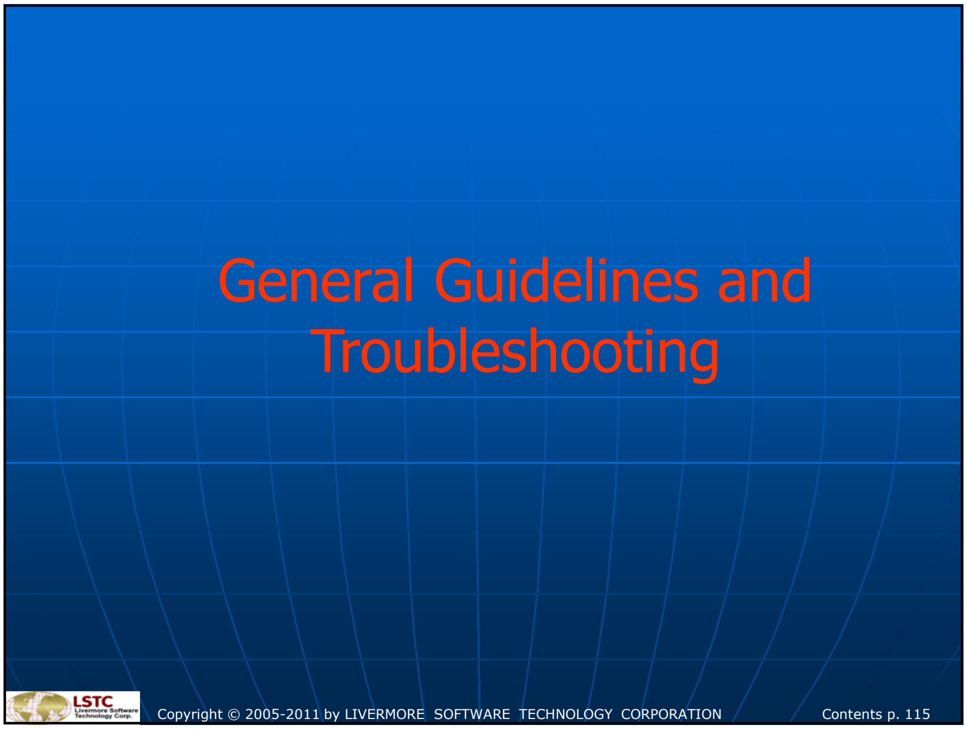
System Tuning

AMD 6200 series

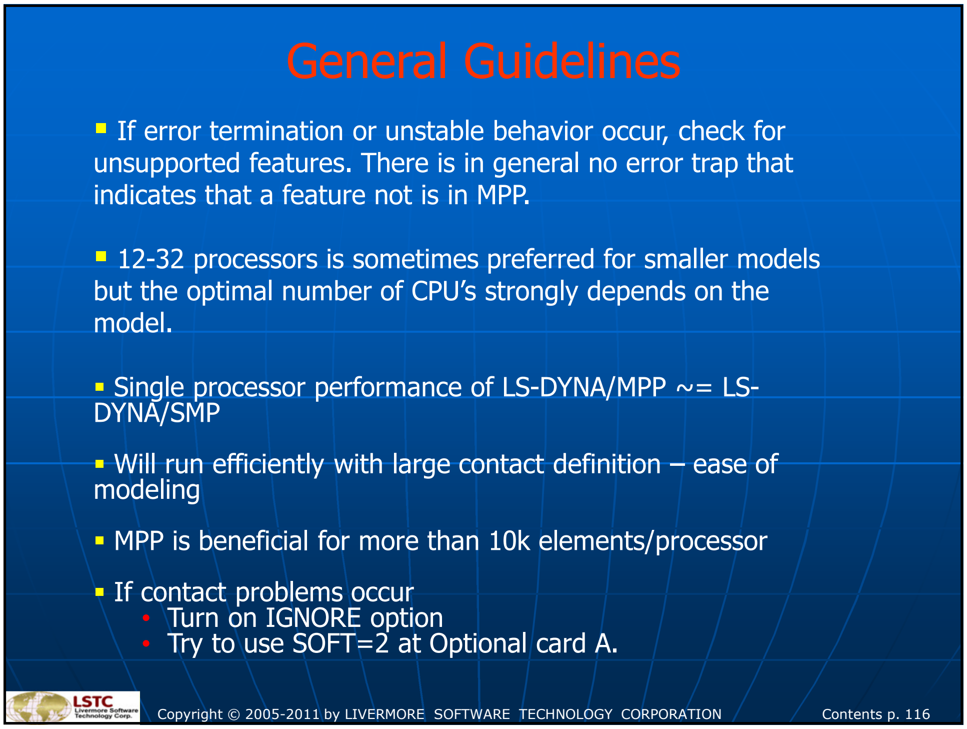
- use only the core has its own FPU and L2 cache under Platform MPI

```
 mpirun -np 4 -cpu_bind=v,MAP_CPU:0,2,4,6
or
 mpirun -np 4 -cpu_bind=v,MAP_CPU:0,8,16,24
```

- Use AMD Open64 AVX enabled exe
  http://ftp.lstc.com/beta/mpp-lsdyna/mpp971/R6.1.0
  (look for exec having the string "Open64avx").

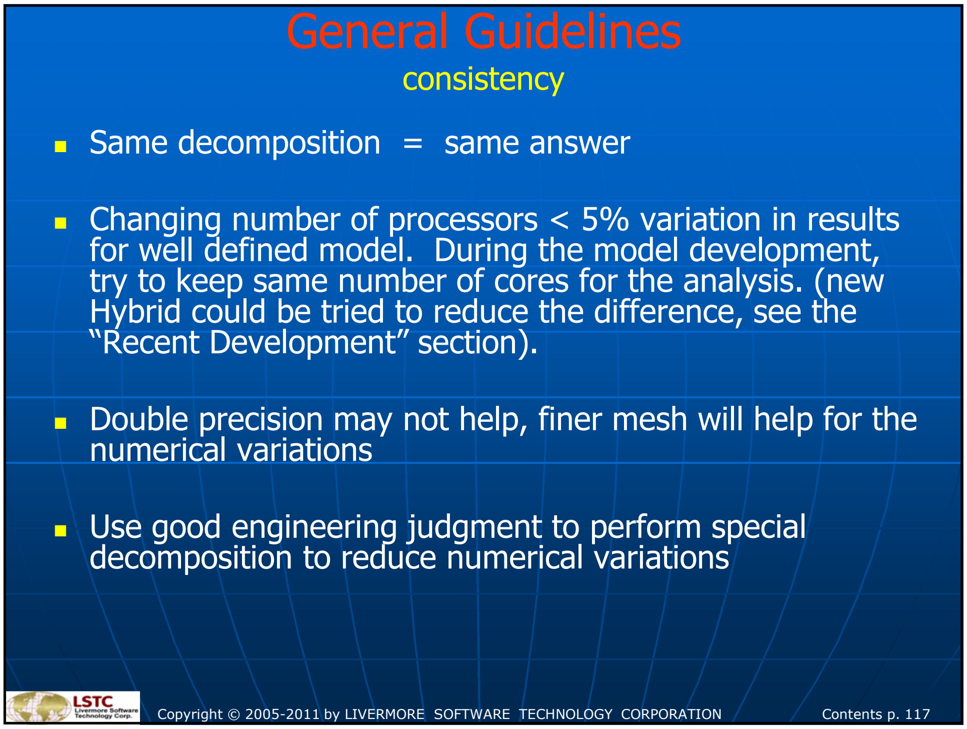


General Guidelines and Troubleshooting



General Guidelines

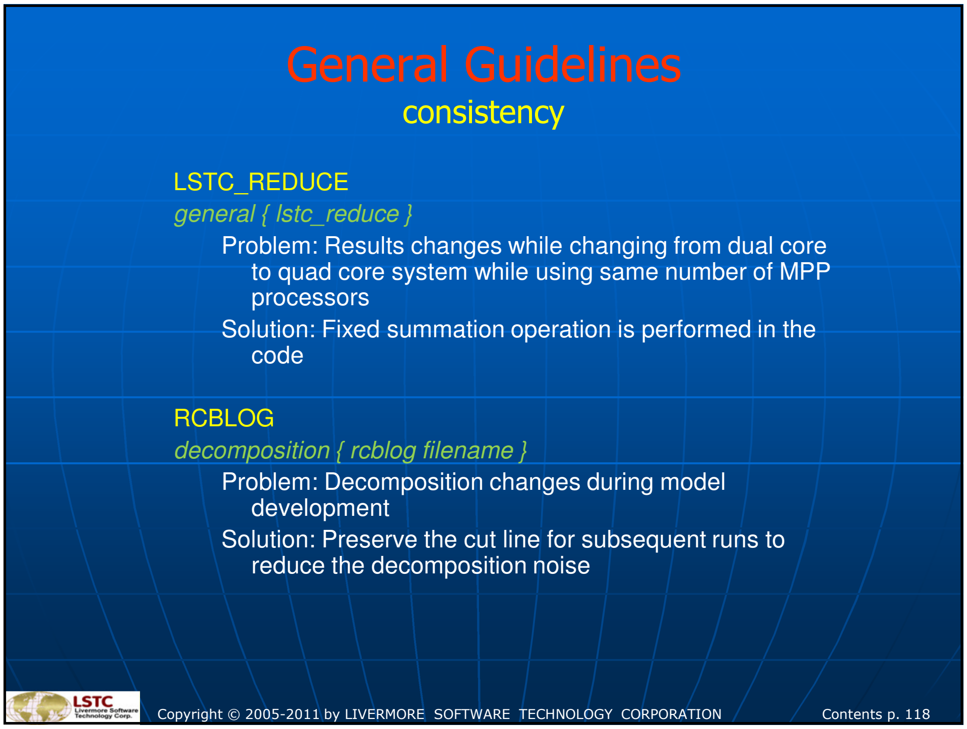
- If error termination or unstable behavior occur, check for unsupported features. There is in general no error trap that indicates that a feature not is in MPP.
- 12-32 processors is sometimes preferred for smaller models but the optimal number of CPU's strongly depends on the model.
- Single processor performance of LS-DYNA/MPP ~= LS-DYNA/SMP
- Will run efficiently with large contact definition – ease of modeling
- MPP is beneficial for more than 10k elements/processor
- If contact problems occur
  - Turn on IGNORE option
  - Try to use SOFT=2 at Optional card A.



General Guidelines

consistency

- Same decomposition = same answer
- Changing number of processors < 5% variation in results for well defined model. During the model development, try to keep same number of cores for the analysis. (new Hybrid could be tried to reduce the difference, see the "Recent Development" section).
- Double precision may not help, finer mesh will help for the numerical variations
- Use good engineering judgment to perform special decomposition to reduce numerical variations



General Guidelines

consistency

LSTC_REDUCE

general { lstc_reduce }

Problem: Results changes while changing from dual core to quad core system while using same number of MPP processors

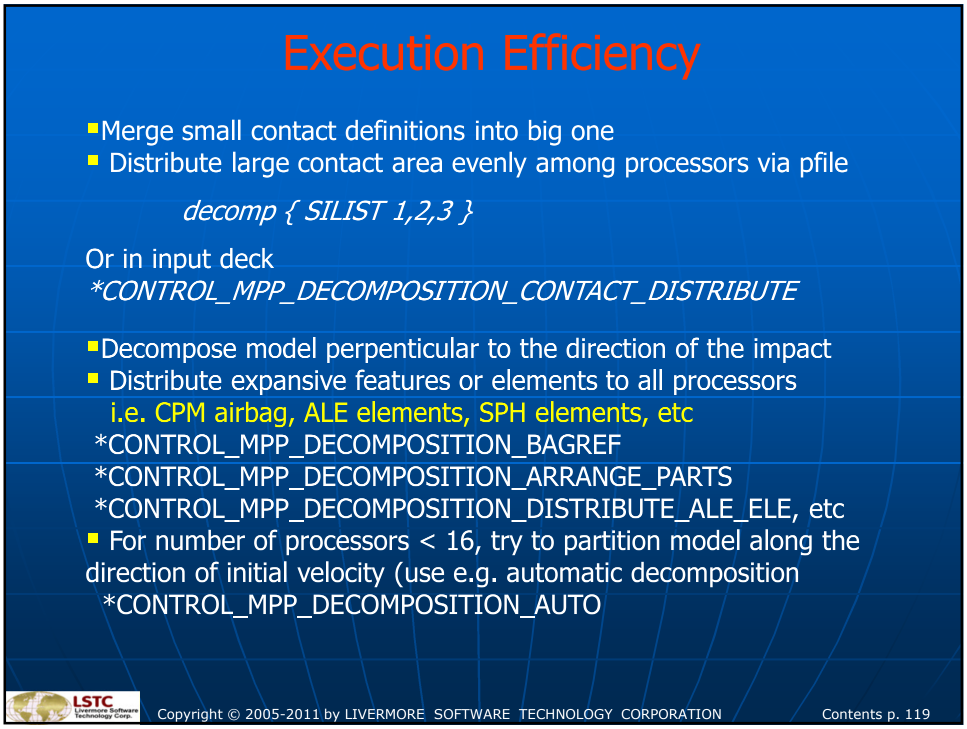
Solution: Fixed summation operation is performed in the code

RCBLOG

decomposition { rcblog filename }

Problem: Decomposition changes during model development

Solution: Preserve the cut line for subsequent runs to reduce the decomposition noise



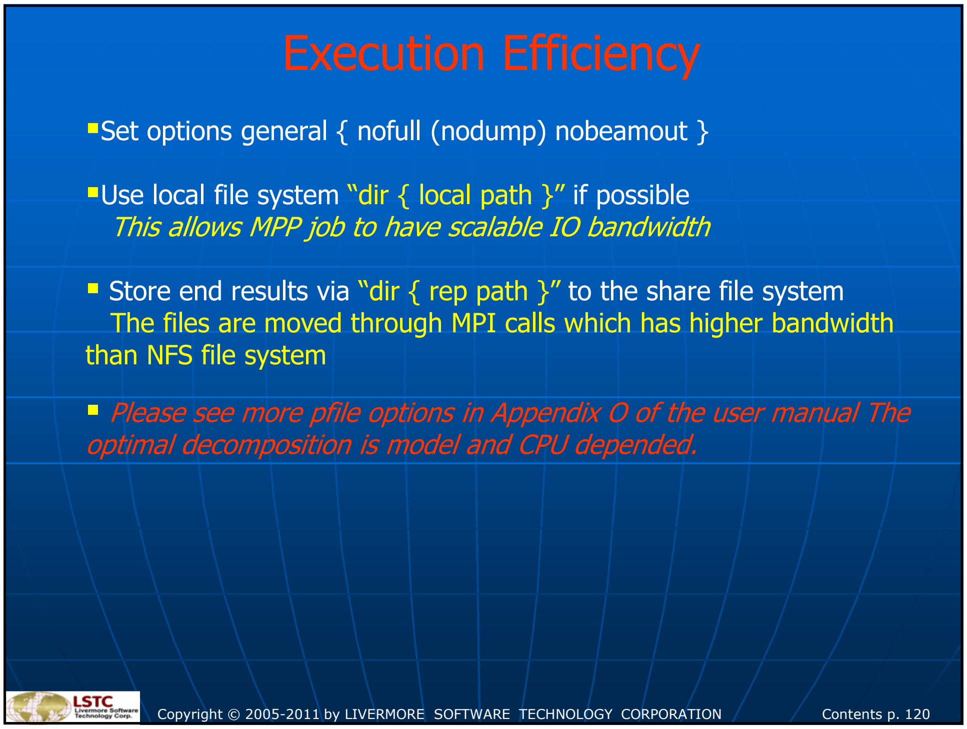
Execution Efficiency

- Merge small contact definitions into big one
- Distribute large contact area evenly among processors via pfile

decomp { SILIST 1,2,3 }

Or in input deck
**CONTROL_MPP_DECOMPOSITION_CONTACT_DISTRIBUTE*

- Decompose model perpenticular to the direction of the impact
- Distribute expansive features or elements to all processors
  i.e. CPM airbag, ALE elements, SPH elements, etc
  *CONTROL_MPP_DECOMPOSITION_BAGREF
  *CONTROL_MPP_DECOMPOSITION_ARRANGE_PARTS
  *CONTROL_MPP_DECOMPOSITION_DISTRIBUTE_ALE_ELE, etc
- For number of processors < 16, try to partition model along the direction of initial velocity (use e.g. automatic decomposition
  *CONTROL_MPP_DECOMPOSITION_AUTO



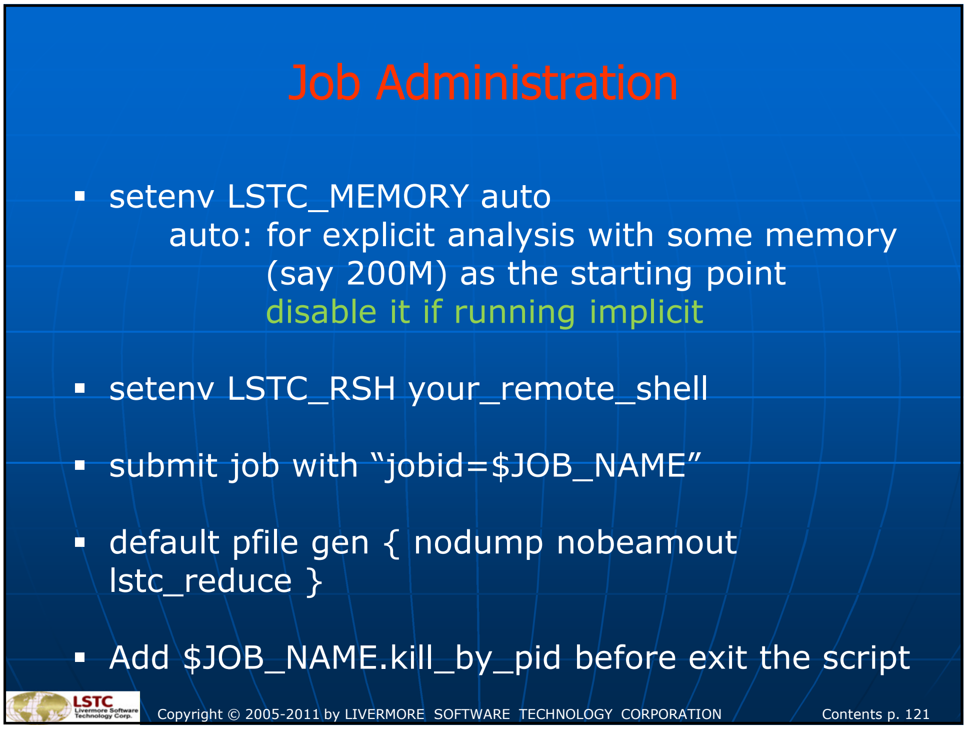
Execution Efficiency

- Set options general { nofull (nodump) nobeamout }

- Use local file system "dir { local path }" if possible
  This allows MPP job to have scalable IO bandwidth

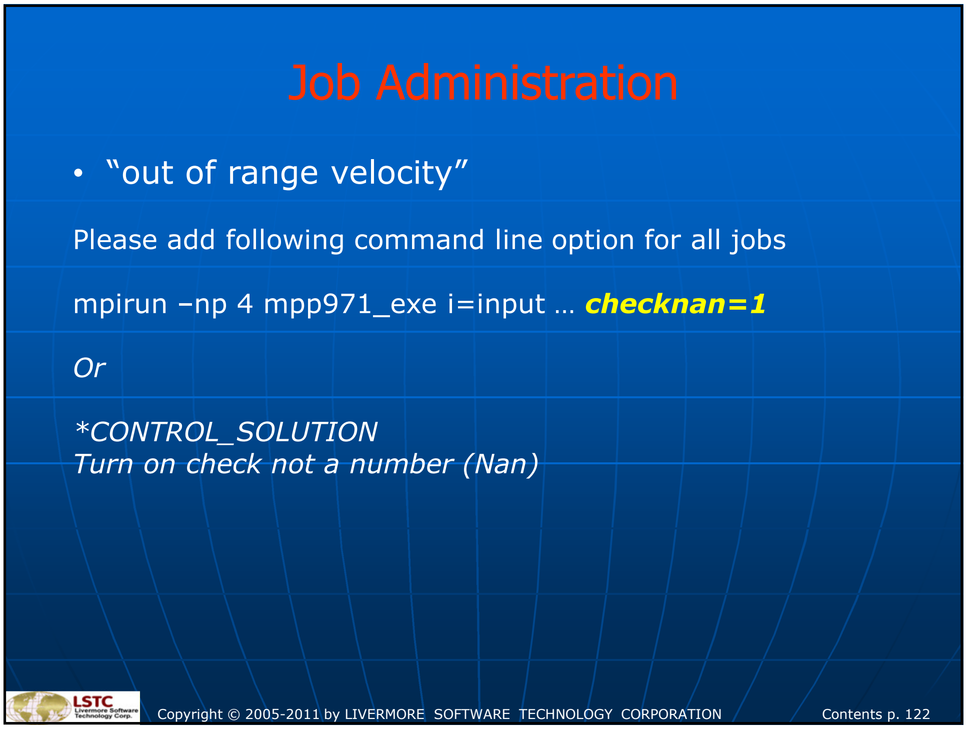
- Store end results via "dir { rep path }" to the share file system
  The files are moved through MPI calls which has higher bandwidth than NFS file system

- *Please see more pfile options in Appendix O of the user manual The optimal decomposition is model and CPU depended.*



Job Administration

- setenv LSTC_MEMORY auto
  auto: for explicit analysis with some memory (say 200M) as the starting point
  disable it if running implicit
- setenv LSTC_RSH your_remote_shell
- submit job with “jobid=$JOB_NAME”
- default pfile gen { nodump nobeamout lstc_reduce }
- Add $JOB_NAME.kill_by_pid before exit the script



Job Administration

- “out of range velocity”

Please add following command line option for all jobs

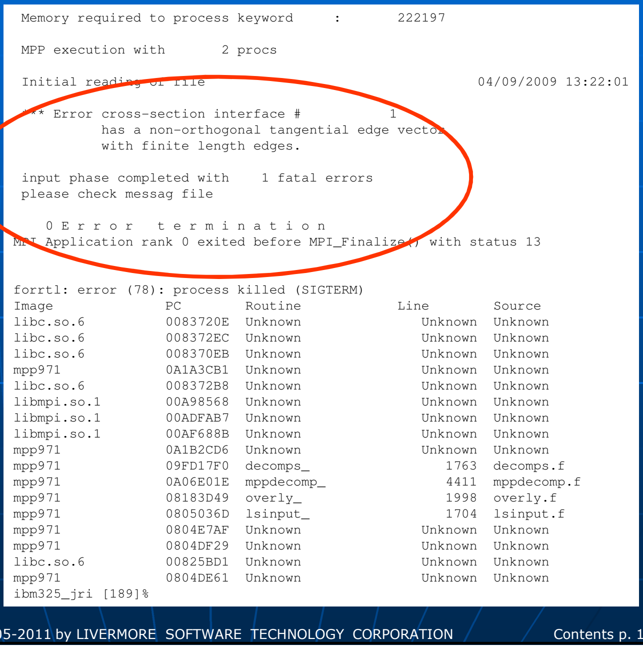
mpirun –np 4 mpp971_exe i=input … ***checknan=1***

Or

**CONTROL_SOLUTION*
Turn on check not a number (Nan)

Troubleshooting

Problem



```
 Memory required to process keyword     :       222197

 MPP execution with       2 procs

 Initial reading of file                                      04/09/2009 13:22:01

 *** Error cross-section interface #            1
          has a non-orthogonal tangential edge vector
          with finite length edges.

 input phase completed with    1 fatal errors
 please check messag file

    0 E r r o r   t e r m i n a t i o n
MPI Application rank 0 exited before MPI_Finalize() with status 13

forrtl: error (78): process killed (SIGTERM)
Image              PC        Routine            Line        Source
libc.so.6          0083720E  Unknown               Unknown  Unknown
libc.so.6          008372EC  Unknown               Unknown  Unknown
libc.so.6          008370EB  Unknown               Unknown  Unknown
mpp971             0A1A3CB1  Unknown               Unknown  Unknown
libc.so.6          008372B8  Unknown               Unknown  Unknown
libmpi.so.1        00A98568  Unknown               Unknown  Unknown
libmpi.so.1        00ADFAB7  Unknown               Unknown  Unknown
libmpi.so.1        00AF688B  Unknown               Unknown  Unknown
mpp971             0A1B2CD6  Unknown               Unknown  Unknown
mpp971             09FD17F0  decomps_                 1763  decomps.f
mpp971             0A06E01E  mppdecomp_               4411  mppdecomp.f
mpp971             08183D49  overly_                  1998  overly.f
mpp971             0805036D  lsinput_                 1704  lsinput.f
mpp971             0804E7AF  Unknown               Unknown  Unknown
mpp971             0804DF29  Unknown               Unknown  Unknown
libc.so.6          00825BD1  Unknown               Unknown  Unknown
mpp971             0804DE61  Unknown               Unknown  Unknown
ibm325_jri [189]%
```

In MPP the error can look serious!

Troubleshooting

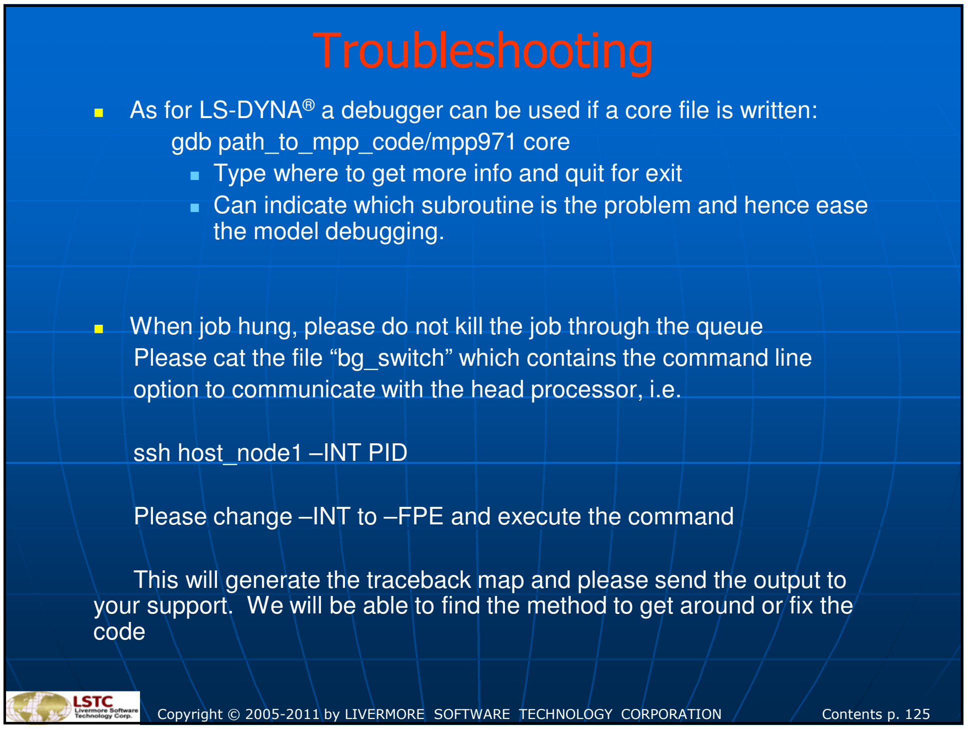
- Find the “nodelist” file which contains the list of hostname used for the current job.
  node001
  node002
  etc

- The error messages from MPP-DYNA can be different from LS-DYNA®

- To locate an error one often has to search each of the messag files mes#### in order to find any information. These files are written for each processor.

- The code will trap the segmentation violation (SEGV) and output the rank number. One could rerun the job and attach the debugger to the running thread and get the trace back map. This usually gives good information for changing input.

```
gdb path_to_mpp_code/mpp971 PID
> continue
SEGV
> where
```



Troubleshooting

- As for LS-DYNA® a debugger can be used if a core file is written:

```
gdb path_to_mpp_code/mpp971 core
```

  - Type where to get more info and quit for exit
  - Can indicate which subroutine is the problem and hence ease the model debugging.

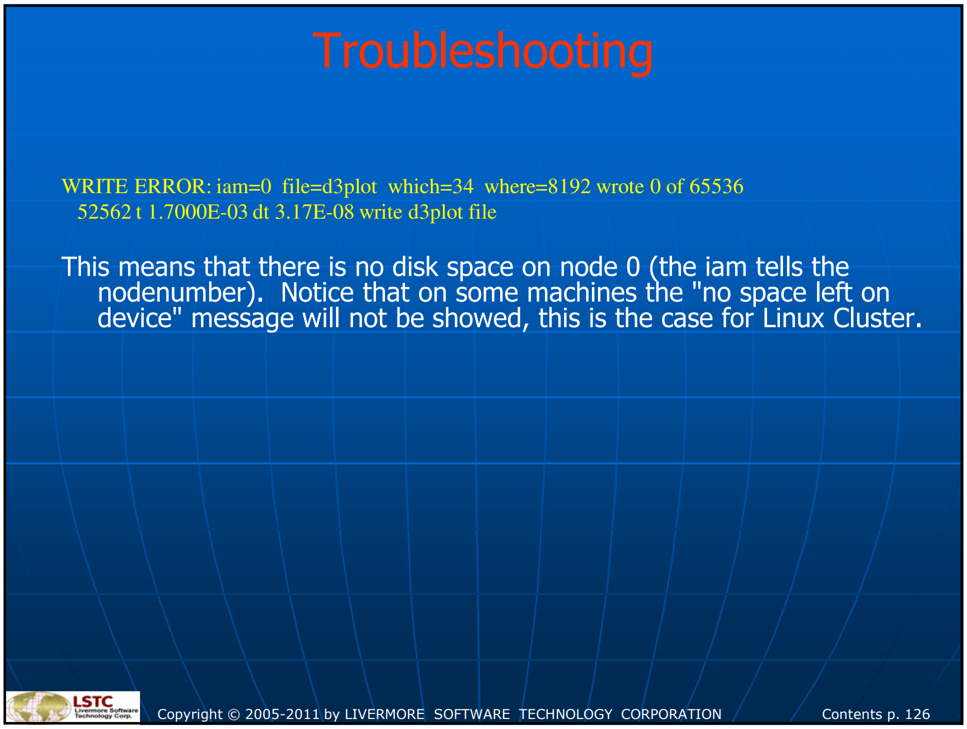
- When job hung, please do not kill the job through the queue

  Please cat the file “bg_switch” which contains the command line option to communicate with the head processor, i.e.

```
ssh host_node1 –INT PID
```

  Please change –INT to –FPE and execute the command

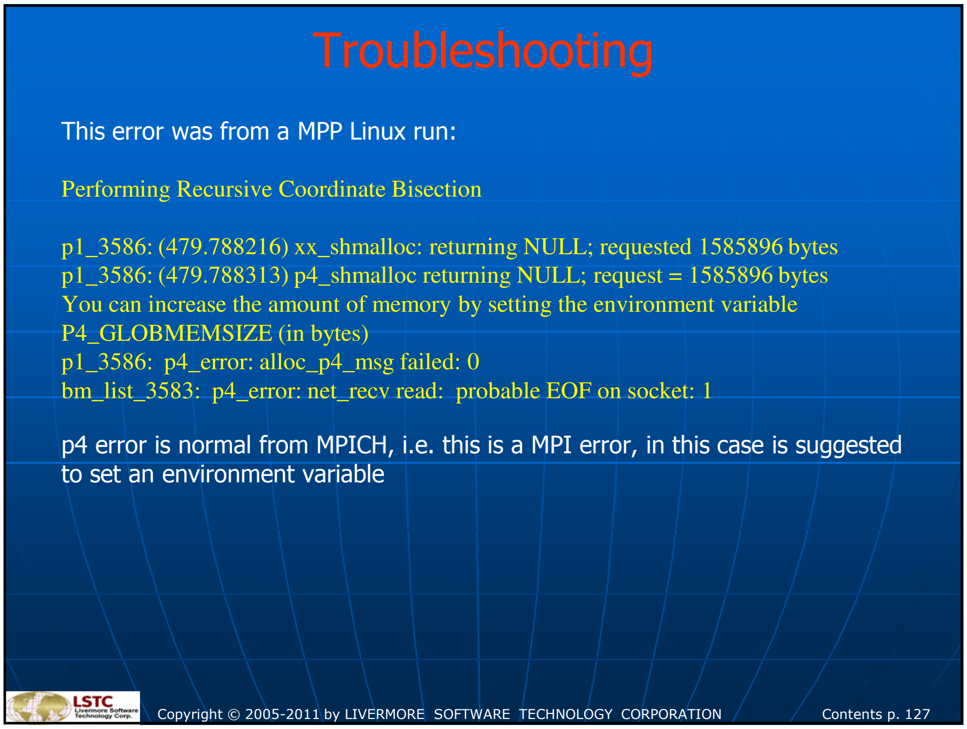
  This will generate the traceback map and please send the output to your support. We will be able to find the method to get around or fix the code



Troubleshooting

```
WRITE ERROR: iam=0  file=d3plot  which=34  where=8192 wrote 0 of 65536
   52562 t 1.7000E-03 dt 3.17E-08 write d3plot file
```

This means that there is no disk space on node 0 (the iam tells the nodenumber). Notice that on some machines the "no space left on device" message will not be showed, this is the case for Linux Cluster.



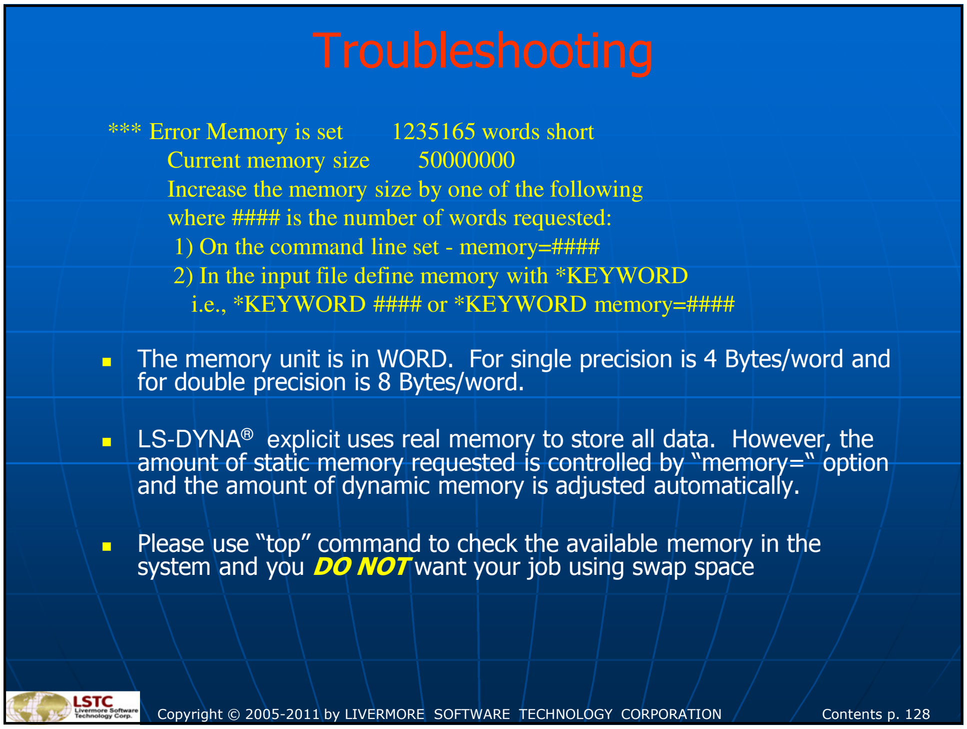
Troubleshooting

This error was from a MPP Linux run:

```
Performing Recursive Coordinate Bisection

p1_3586: (479.788216) xx_shmalloc: returning NULL; requested 1585896 bytes
p1_3586: (479.788313) p4_shmalloc returning NULL; request = 1585896 bytes
You can increase the amount of memory by setting the environment variable
P4_GLOBMEMSIZE (in bytes)
p1_3586:  p4_error: alloc_p4_msg failed: 0
bm_list_3583:  p4_error: net_recv read:  probable EOF on socket: 1
```

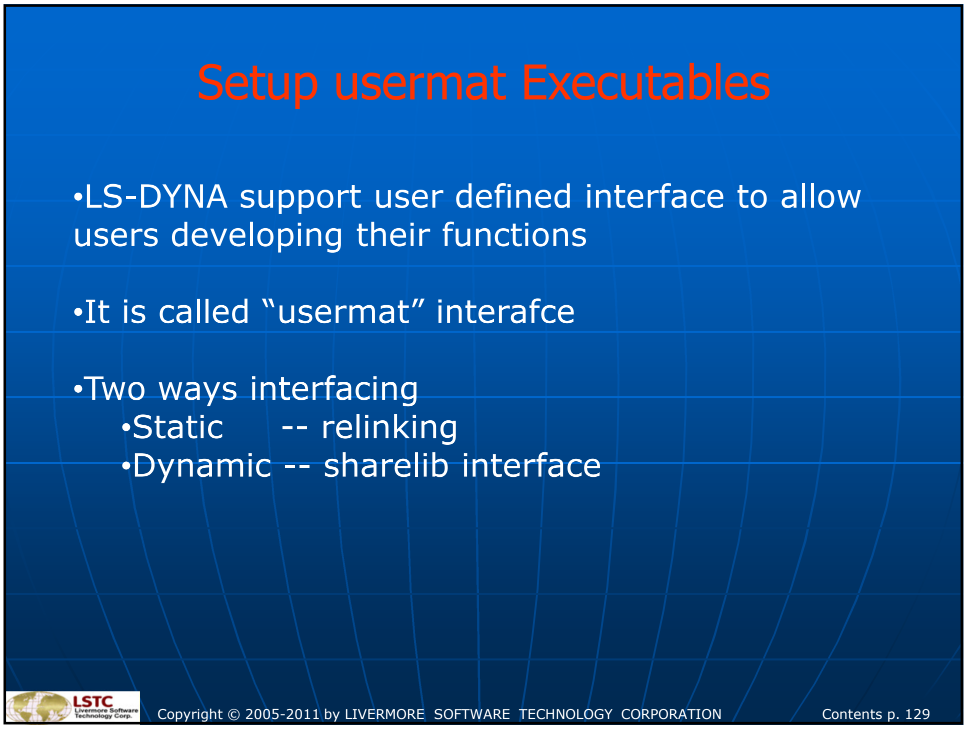
p4 error is normal from MPICH, i.e. this is a MPI error, in this case is suggested to set an environment variable



Troubleshooting

```
*** Error Memory is set        1235165 words short
        Current memory size        50000000
        Increase the memory size by one of the following
        where #### is the number of words requested:
         1) On the command line set - memory=####
         2) In the input file define memory with *KEYWORD
            i.e., *KEYWORD #### or *KEYWORD memory=####
```

- The memory unit is in WORD. For single precision is 4 Bytes/word and for double precision is 8 Bytes/word.
- LS-DYNA® explicit uses real memory to store all data. However, the amount of static memory requested is controlled by "memory=" option and the amount of dynamic memory is adjusted automatically.
- Please use "top" command to check the available memory in the system and you ***DO NOT*** want your job using swap space

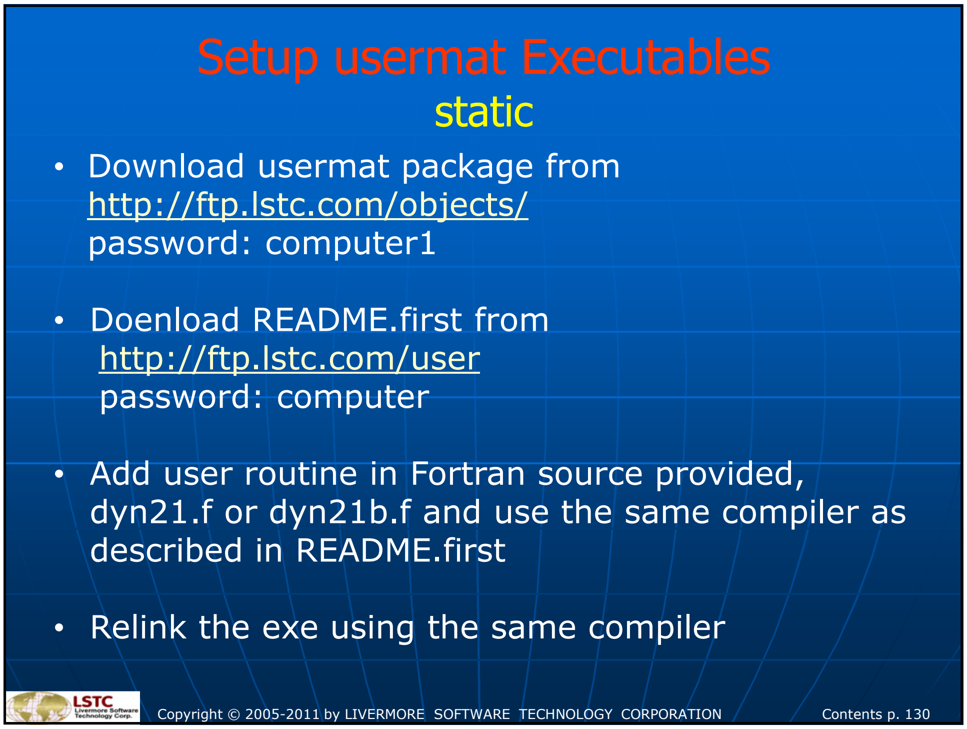


Setup usermat Executables

•LS-DYNA support user defined interface to allow users developing their functions

•It is called “usermat” interafce

•Two ways interfacing
  - •Static -- relinking
  - •Dynamic -- sharelib interface



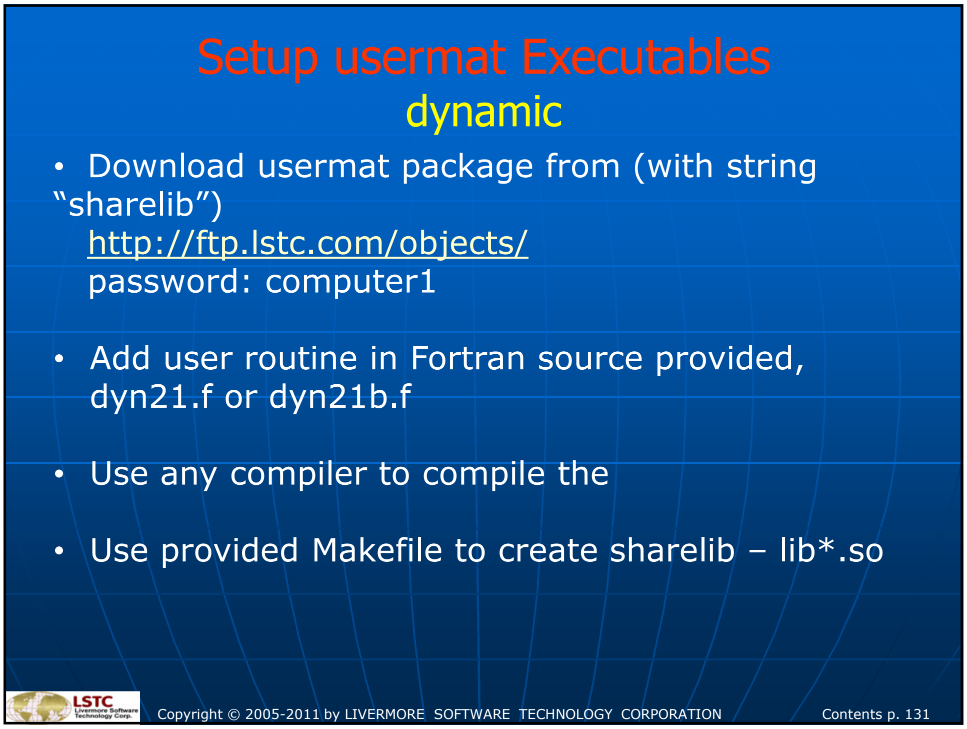
Setup usermat Executables
static

- Download usermat package from
  http://ftp.lstc.com/objects/
  password: computer1

- Doenload README.first from
  http://ftp.lstc.com/user
  password: computer

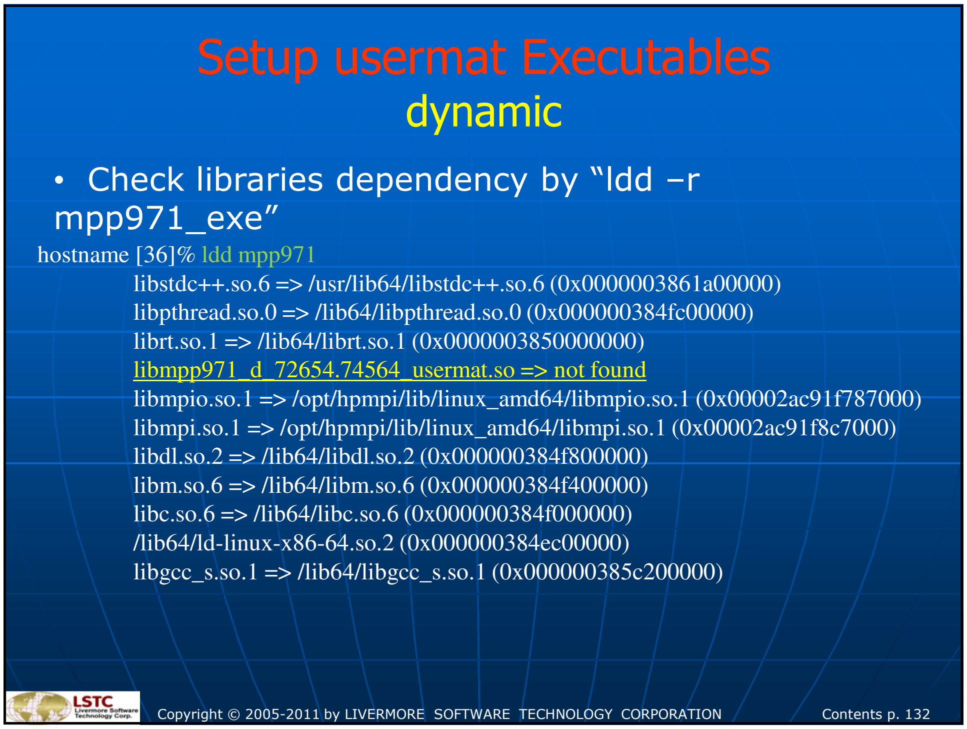
- Add user routine in Fortran source provided, dyn21.f or dyn21b.f and use the same compiler as described in README.first

- Relink the exe using the same compiler



Setup usermat Executables
dynamic

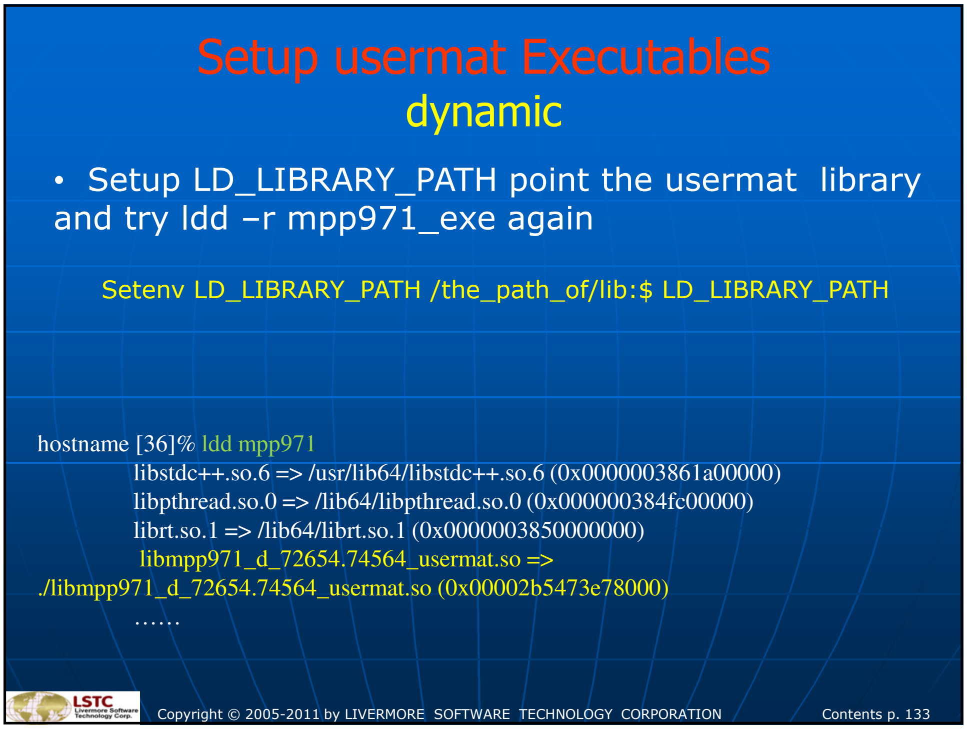
- Download usermat package from (with string "sharelib")
  http://ftp.lstc.com/objects/
  password: computer1
- Add user routine in Fortran source provided, dyn21.f or dyn21b.f
- Use any compiler to compile the
- Use provided Makefile to create sharelib – lib*.so



Setup usermat Executables
dynamic

- Check libraries dependency by "ldd –r mpp971_exe"

```
hostname [36]% ldd mpp971
        libstdc++.so.6 => /usr/lib64/libstdc++.so.6 (0x0000003861a00000)
        libpthread.so.0 => /lib64/libpthread.so.0 (0x000000384fc00000)
        librt.so.1 => /lib64/librt.so.1 (0x0000003850000000)
        libmpp971_d_72654.74564_usermat.so => not found
        libmpio.so.1 => /opt/hpmpi/lib/linux_amd64/libmpio.so.1 (0x00002ac91f787000)
        libmpi.so.1 => /opt/hpmpi/lib/linux_amd64/libmpi.so.1 (0x00002ac91f8c7000)
        libdl.so.2 => /lib64/libdl.so.2 (0x000000384f800000)
        libm.so.6 => /lib64/libm.so.6 (0x000000384f400000)
        libc.so.6 => /lib64/libc.so.6 (0x000000384f000000)
        /lib64/ld-linux-x86-64.so.2 (0x000000384ec00000)
        libgcc_s.so.1 => /lib64/libgcc_s.so.1 (0x000000385c200000)
```

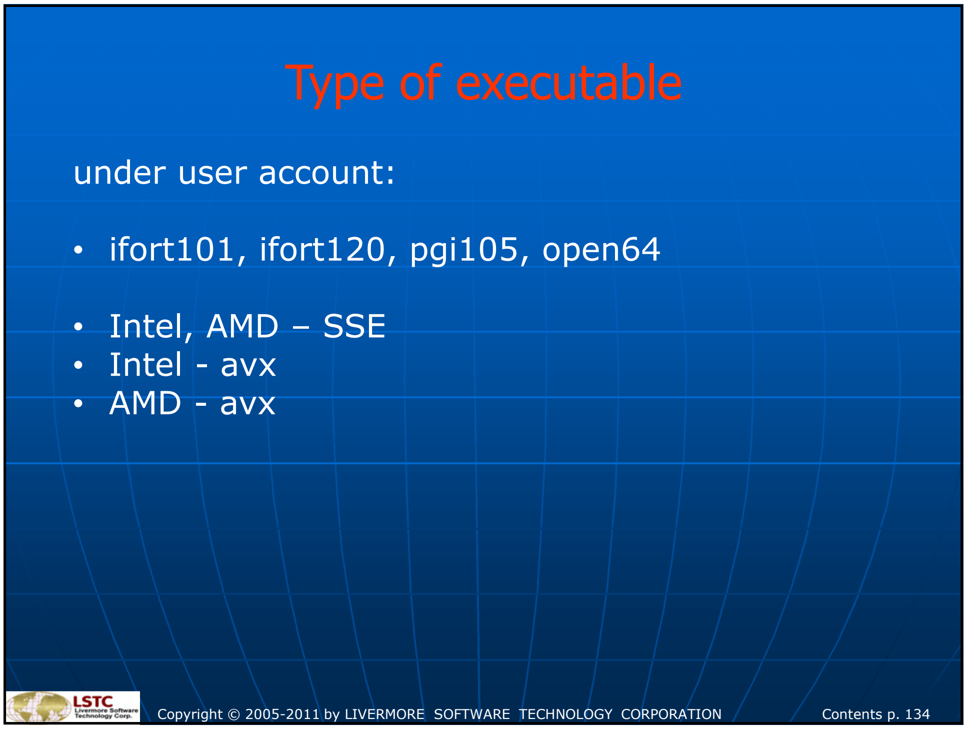


Setup usermat Executables
dynamic

- Setup LD_LIBRARY_PATH point the usermat library and try ldd –r mpp971_exe again

```
Setenv LD_LIBRARY_PATH /the_path_of/lib:$ LD_LIBRARY_PATH
```

```
hostname [36]% ldd mpp971
             libstdc++.so.6 => /usr/lib64/libstdc++.so.6 (0x0000003861a00000)
             libpthread.so.0 => /lib64/libpthread.so.0 (0x000000384fc00000)
             librt.so.1 => /lib64/librt.so.1 (0x0000003850000000)
              libmpp971_d_72654.74564_usermat.so =>
./libmpp971_d_72654.74564_usermat.so (0x00002b5473e78000)
             ……
```



Type of executable

under user account:

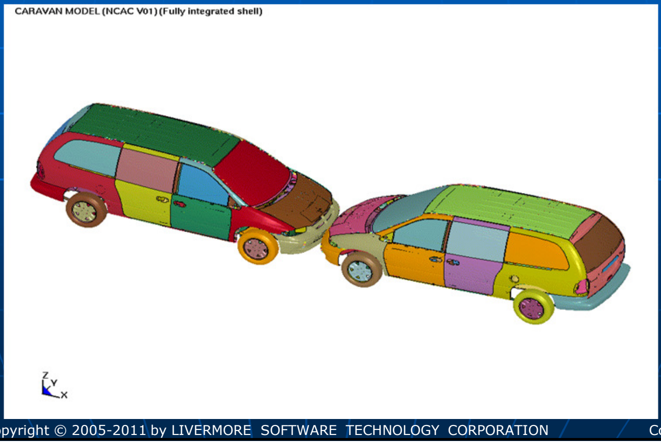
- ifort101, ifort120, pgi105, open64

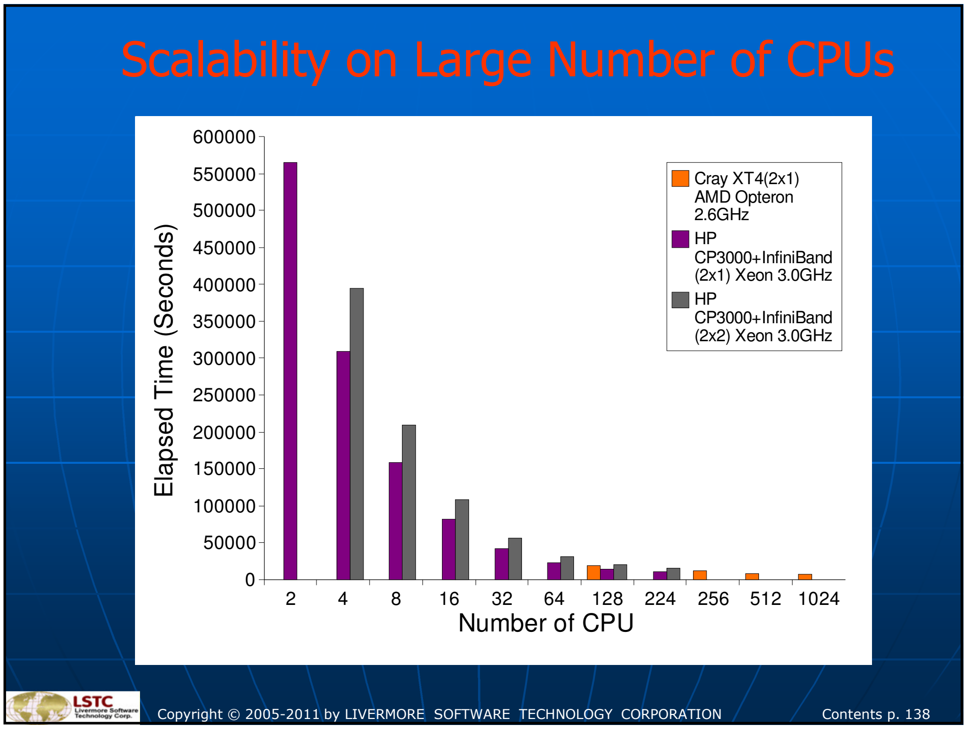
- Intel, AMD – SSE
- Intel - avx
- AMD - avx

Recent Development

Scalability on Large Number of CPUs

Model statistic (car2car model)
~2,500,000 nodes and elements
53 contacts
Fully integrated (type 16) shells





Scalability on Large Number of CPUs

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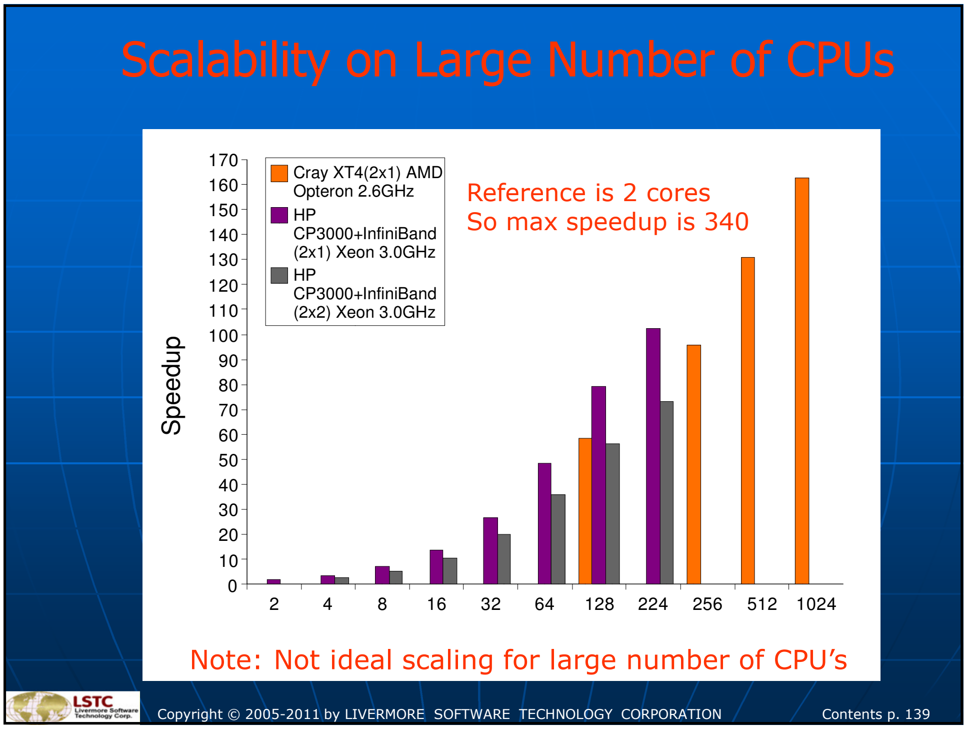

Scalability on Large Number of CPUs

Reference is 2 cores
So max speedup is 340

Note: Not ideal scaling for large number of CPU's

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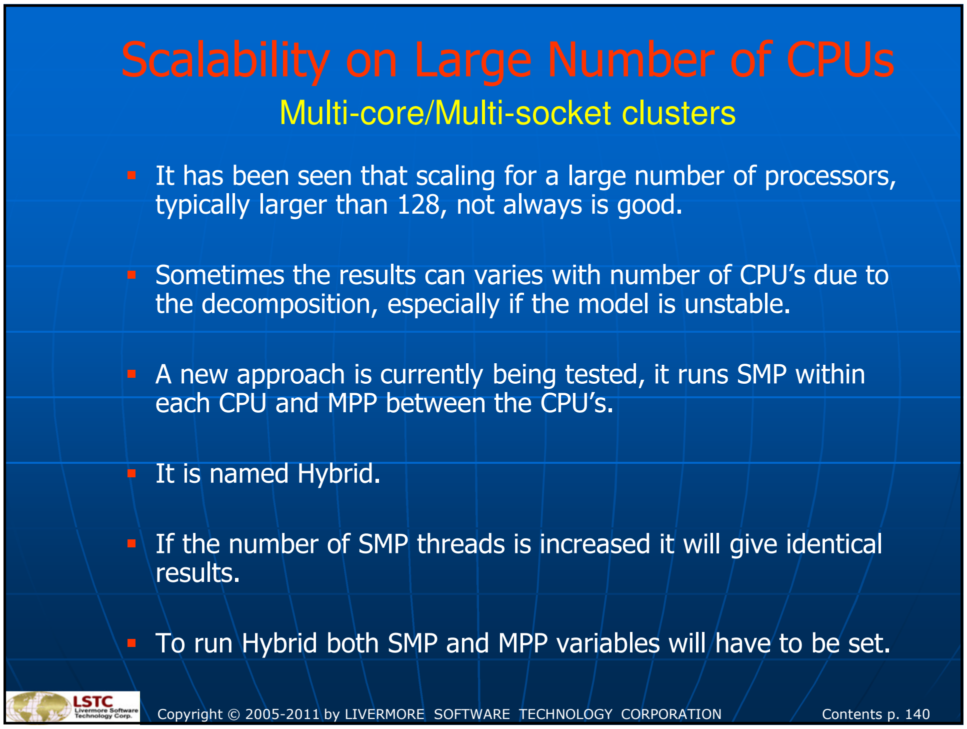
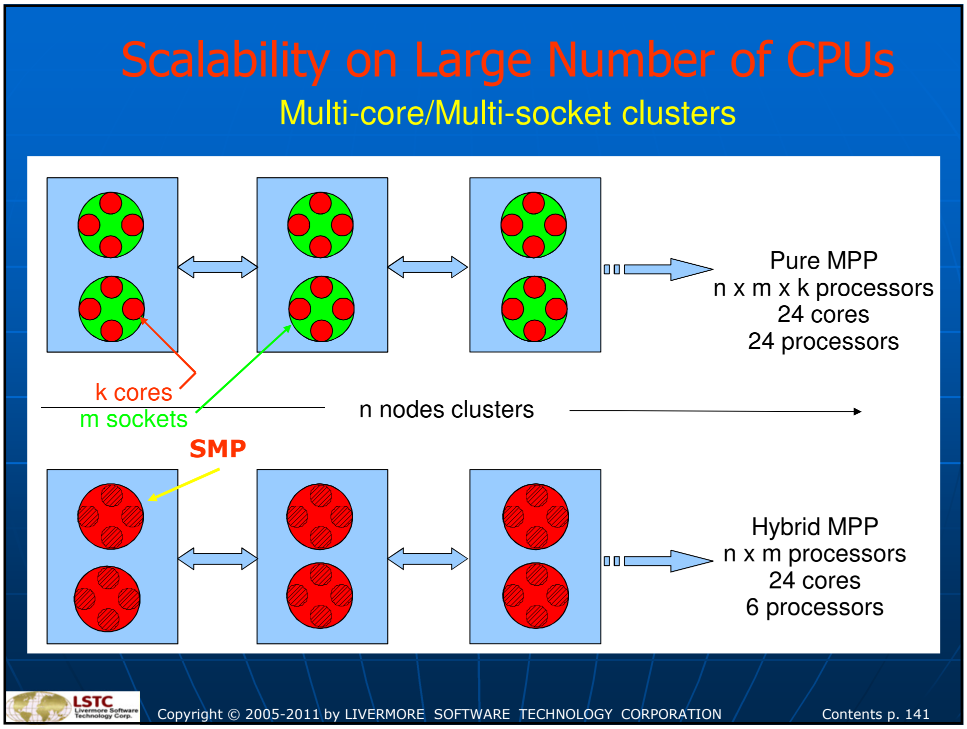


Scalability on Large Number of CPUs

Multi-core/Multi-socket clusters

- It has been seen that scaling for a large number of processors, typically larger than 128, not always is good.
- Sometimes the results can varies with number of CPU's due to the decomposition, especially if the model is unstable.
- A new approach is currently being tested, it runs SMP within each CPU and MPP between the CPU's.
- It is named Hybrid.
- If the number of SMP threads is increased it will give identical results.
- To run Hybrid both SMP and MPP variables will have to be set.



Scalability on Large Number of CPUs

Multi-core/Multi-socket clusters

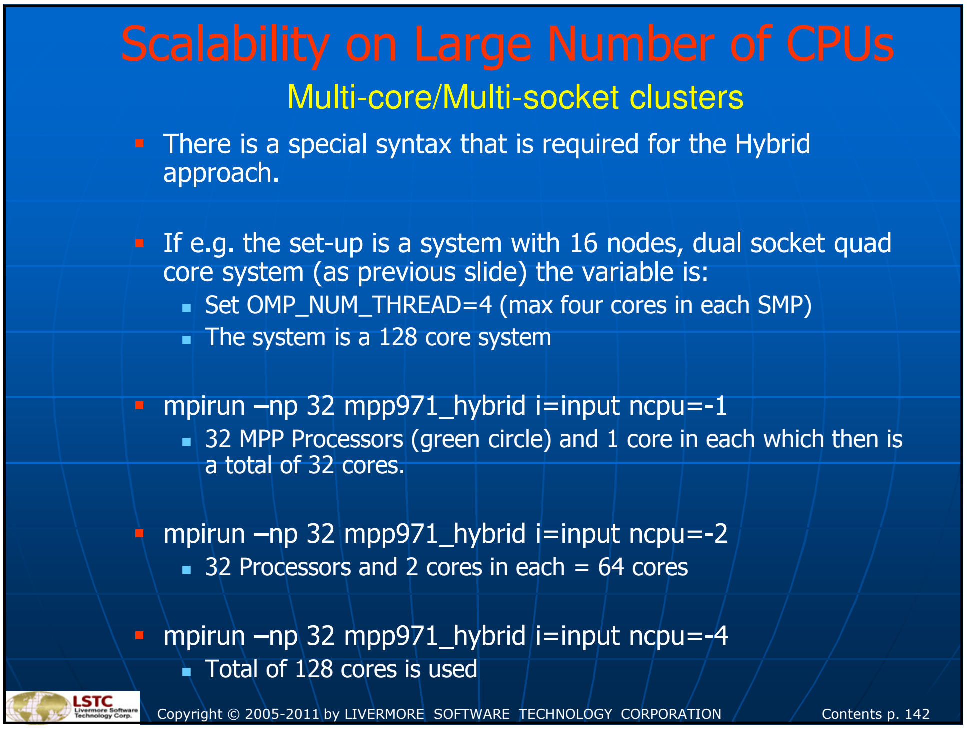
Pure MPP
n x m x k processors
24 cores
24 processors

k cores
m sockets

n nodes clusters

SMP

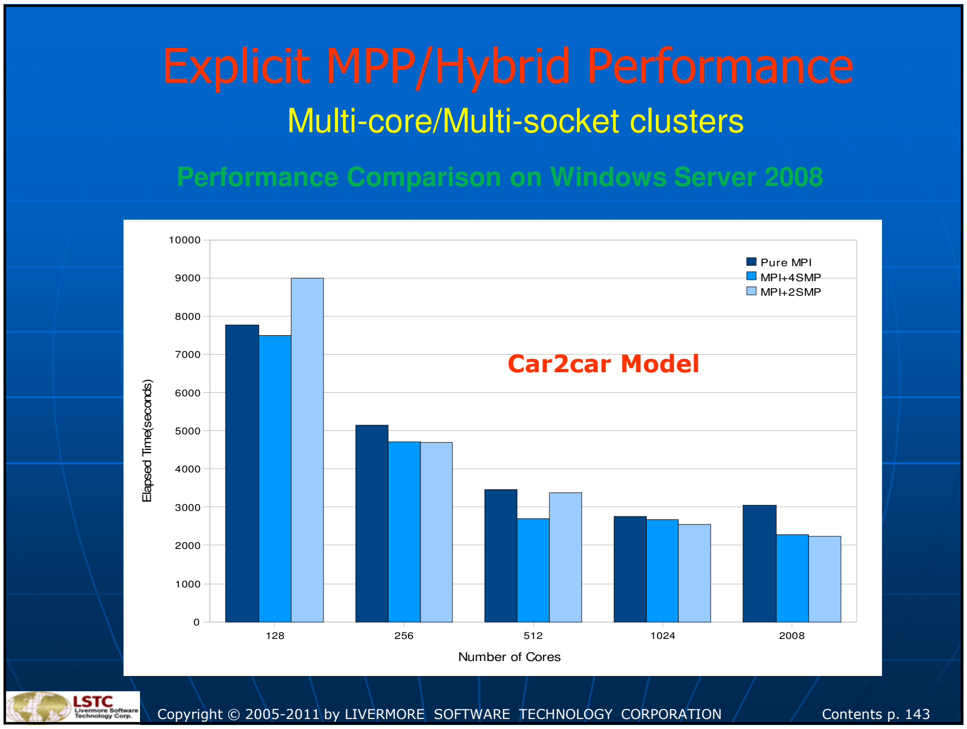
Hybrid MPP
n x m processors
24 cores
6 processors



Scalability on Large Number of CPUs

Multi-core/Multi-socket clusters

- There is a special syntax that is required for the Hybrid approach.
- If e.g. the set-up is a system with 16 nodes, dual socket quad core system (as previous slide) the variable is:
  - Set OMP_NUM_THREAD=4 (max four cores in each SMP)
  - The system is a 128 core system
- mpirun –np 32 mpp971_hybrid i=input ncpu=-1
  - 32 MPP Processors (green circle) and 1 core in each which then is a total of 32 cores.
- mpirun –np 32 mpp971_hybrid i=input ncpu=-2
  - 32 Processors and 2 cores in each = 64 cores
- mpirun –np 32 mpp971_hybrid i=input ncpu=-4
  - Total of 128 cores is used



Explicit MPP/Hybrid Performance

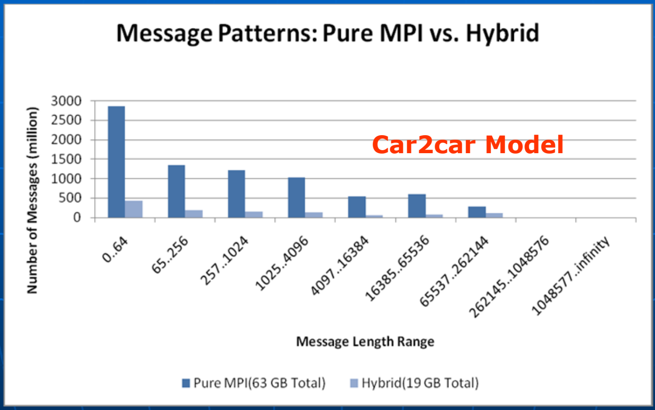
Multi-core/Multi-socket clusters

Performance Comparison on Windows Server 2008

Scalability on Large Number of CPUs

Multi-core/Multi-socket clusters

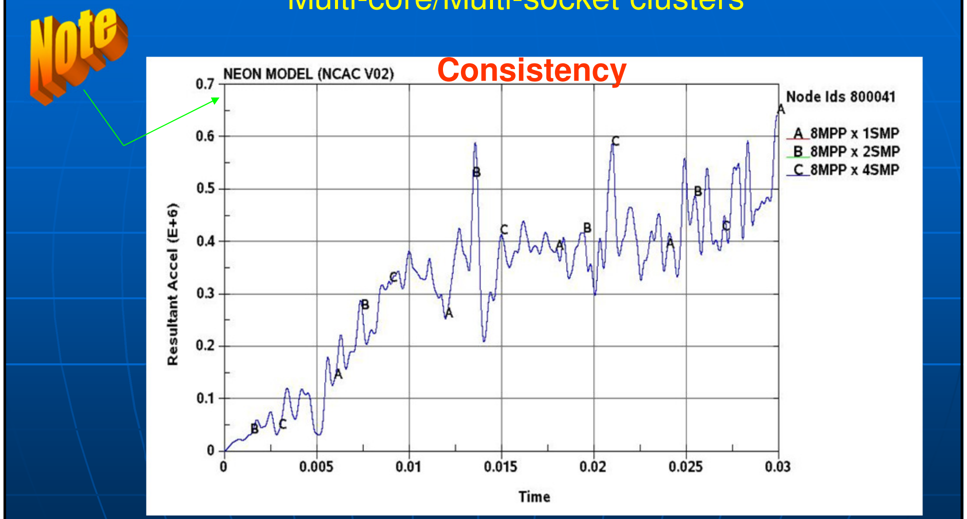
Message Across Network



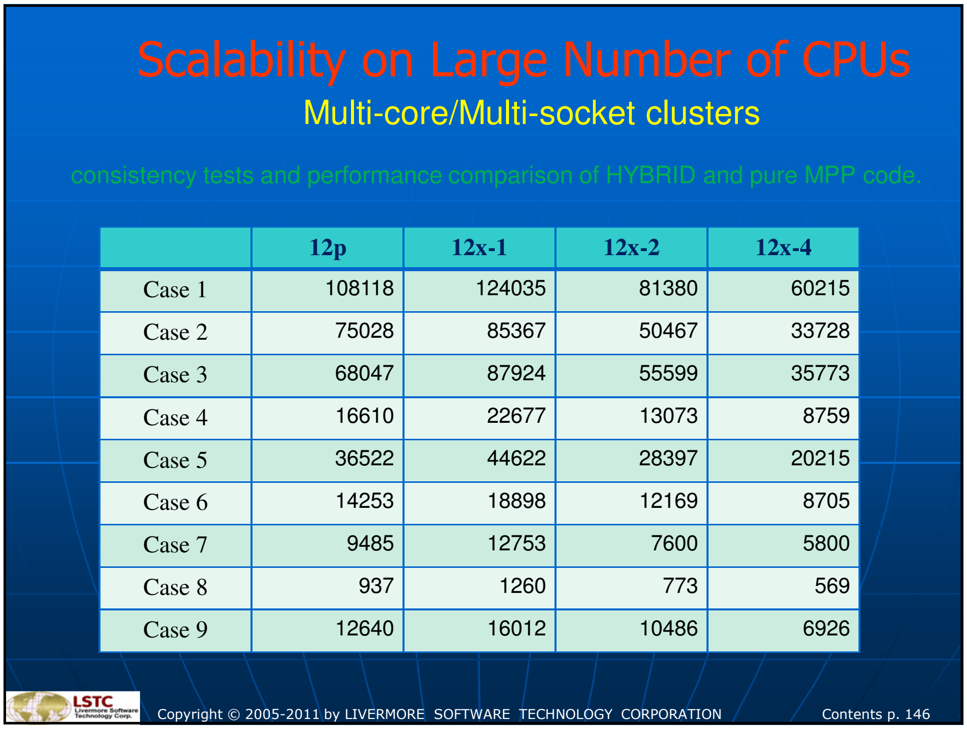
- Hybrid greatly reduce the amount of data through network and provide better scaling to large number of processors

Scalability on Large Number of CPUs

Multi-core/Multi-socket clusters



- Consistent results is be obtained with fix decomposition and changing number of SMP threads

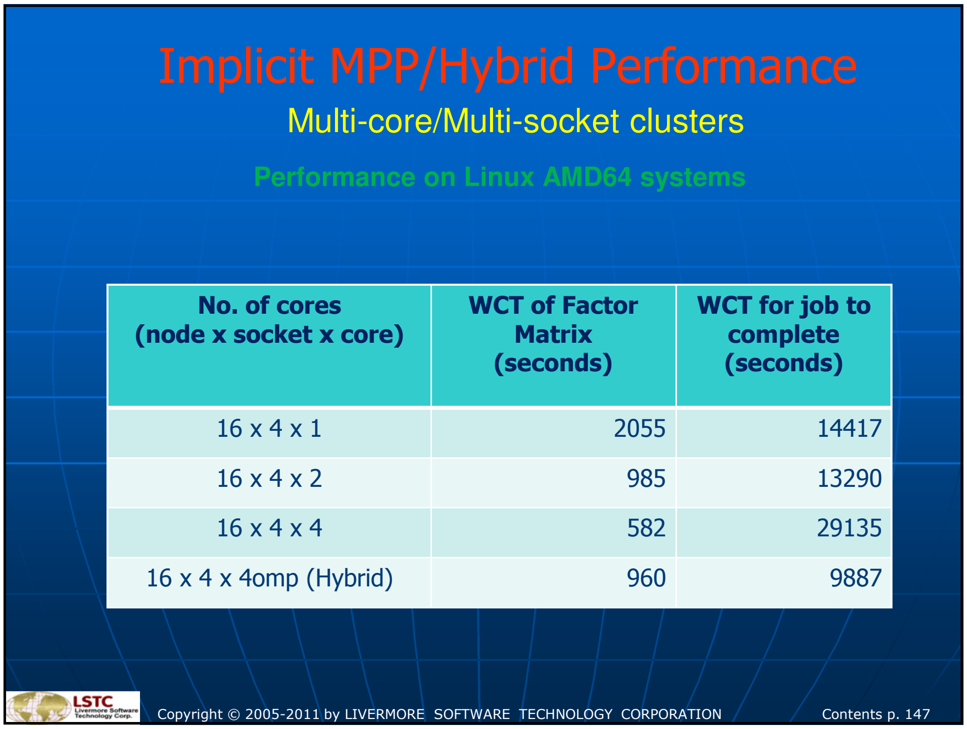


Scalability on Large Number of CPUs

Multi-core/Multi-socket clusters

consistency tests and performance comparison of HYBRID and pure MPP code.

	12p	12x-1	12x-2	12x-4
Case 1	108118	124035	81380	60215
Case 2	75028	85367	50467	33728
Case 3	68047	87924	55599	35773
Case 4	16610	22677	13073	8759
Case 5	36522	44622	28397	20215
Case 6	14253	18898	12169	8705
Case 7	9485	12753	7600	5800
Case 8	937	1260	773	569
Case 9	12640	16012	10486	6926

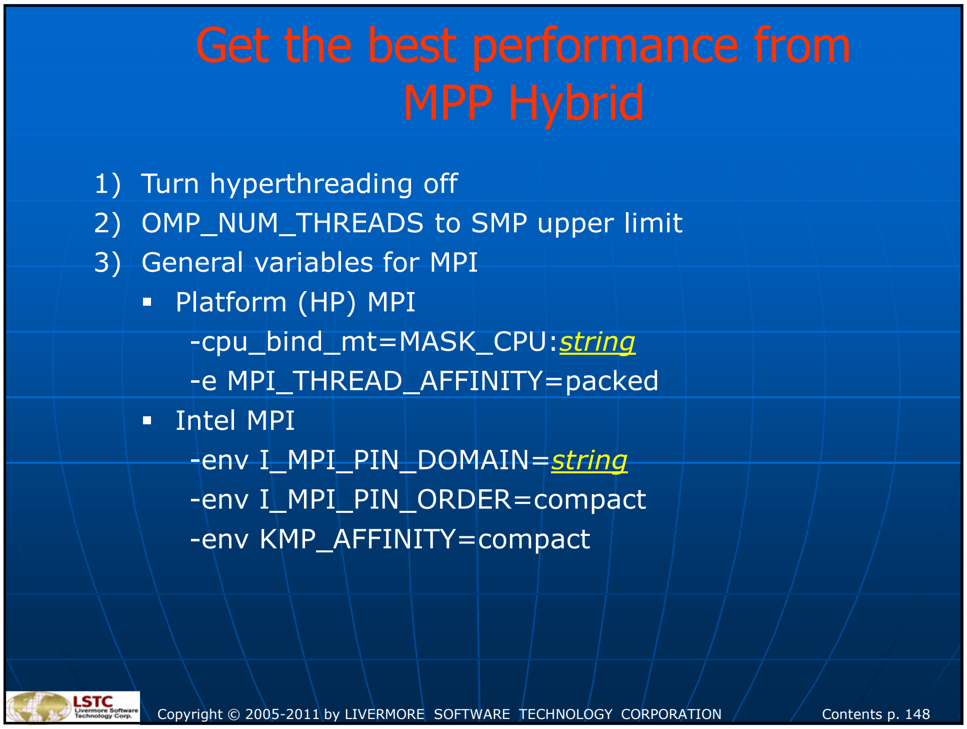


Implicit MPP/Hybrid Performance

Multi-core/Multi-socket clusters

Performance on Linux AMD64 systems

No. of cores (node x socket x core)	WCT of Factor Matrix (seconds)	WCT for job to complete (seconds)
16 x 4 x 1	2055	14417
16 x 4 x 2	985	13290
16 x 4 x 4	582	29135
16 x 4 x 4omp (Hybrid)	960	9887



Get the best performance from MPP Hybrid

1) Turn hyperthreading off
2) OMP_NUM_THREADS to SMP upper limit
3) General variables for MPI
   - Platform (HP) MPI

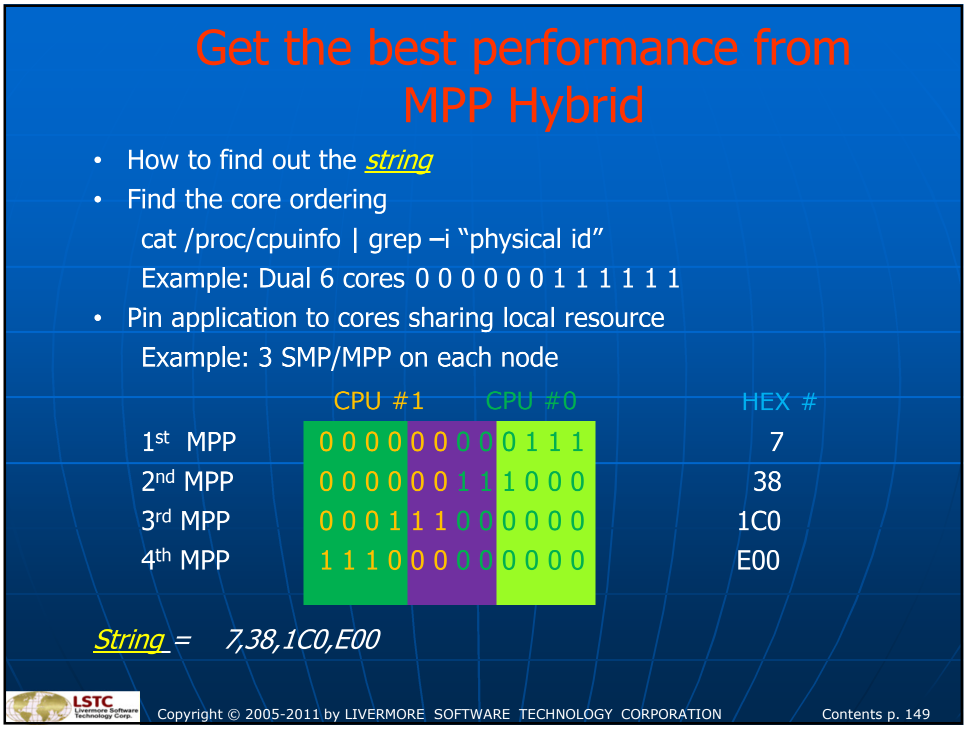
     -cpu_bind_mt=MASK_CPU:*string*

     -e MPI_THREAD_AFFINITY=packed
   - Intel MPI

     -env I_MPI_PIN_DOMAIN=*string*

     -env I_MPI_PIN_ORDER=compact

     -env KMP_AFFINITY=compact



Get the best performance from MPP Hybrid

- How to find out the *string*
- Find the core ordering

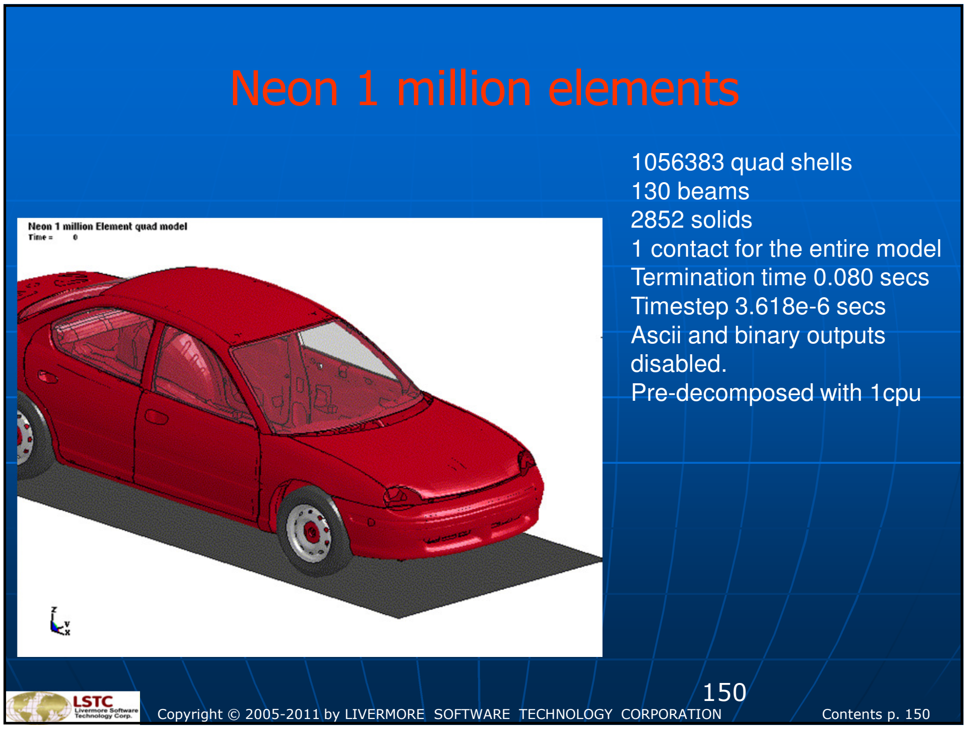
  cat /proc/cpuinfo | grep –i "physical id"

  Example: Dual 6 cores 0 0 0 0 0 0 1 1 1 1 1 1
- Pin application to cores sharing local resource

  Example: 3 SMP/MPP on each node

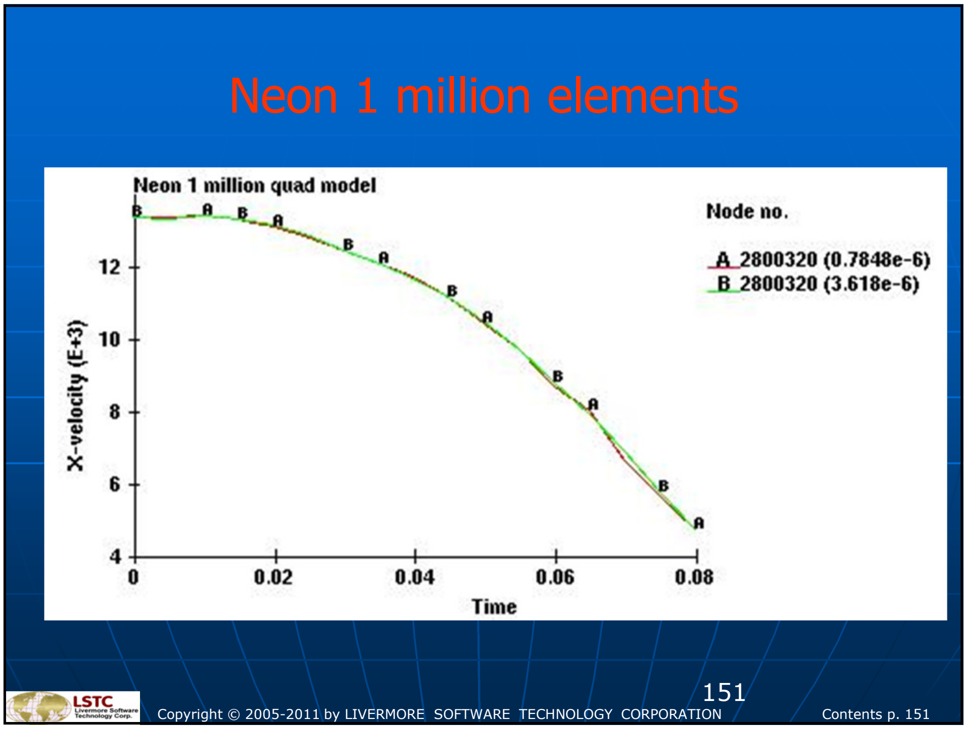
	CPU #1	CPU #0	HEX #
1st MPP	0 0 0 0 0 0	0 0 0 1 1 1	7
2nd MPP	0 0 0 0 0 0	1 1 1 0 0 0	38
3rd MPP	0 0 0 1 1 1	0 0 0 0 0 0	1C0
4th MPP	1 1 1 0 0 0	0 0 0 0 0 0	E00

String = *7,38,1C0,E00*

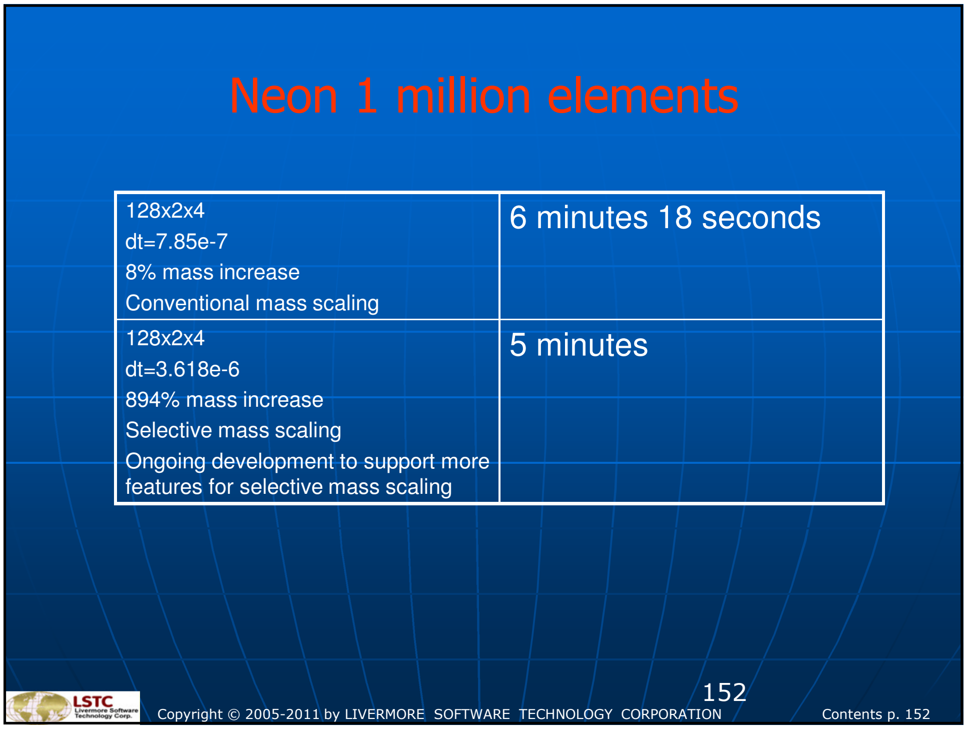


Neon 1 million elements

1056383 quad shells
130 beams
2852 solids
1 contact for the entire model
Termination time 0.080 secs
Timestep 3.618e-6 secs
Ascii and binary outputs disabled.
Pre-decomposed with 1cpu

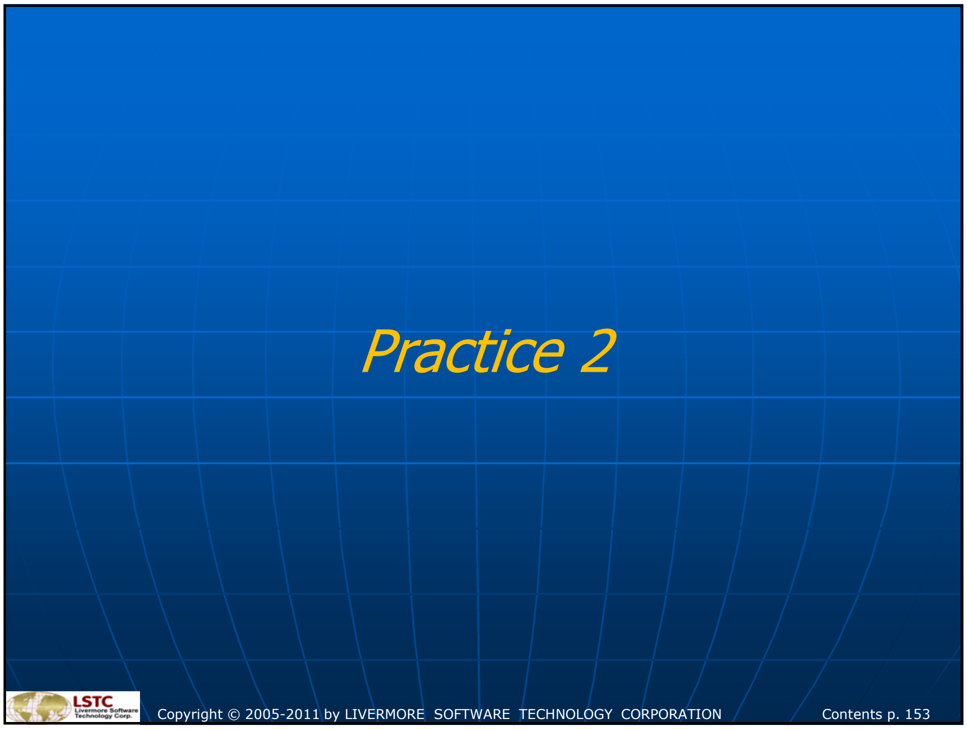


Neon 1 million elements

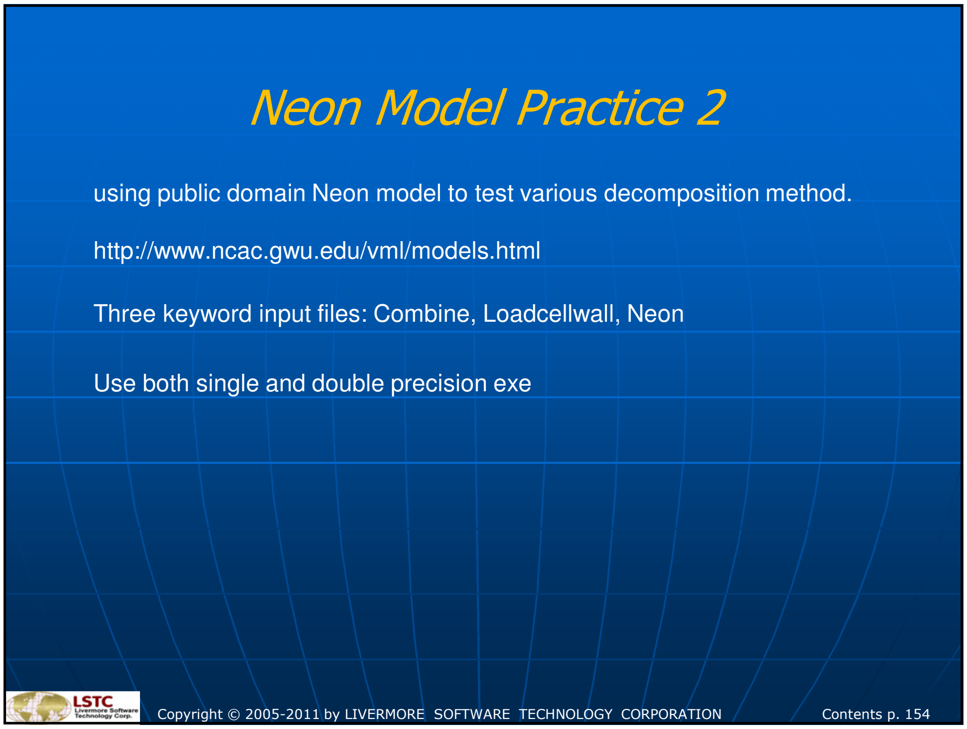


Neon 1 million elements

128x2x4 dt=7.85e-7 8% mass increase Conventional mass scaling	6 minutes 18 seconds
128x2x4 dt=3.618e-6 894% mass increase Selective mass scaling Ongoing development to support more features for selective mass scaling	5 minutes



Practice 2



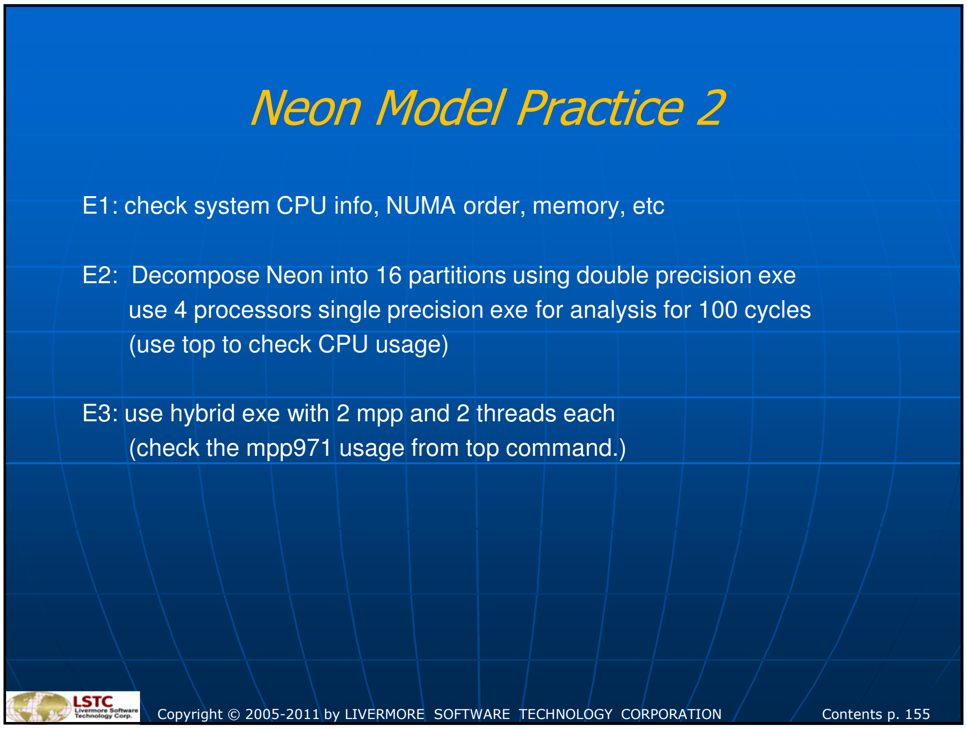
Neon Model Practice 2

using public domain Neon model to test various decomposition method.

http://www.ncac.gwu.edu/vml/models.html

Three keyword input files: Combine, Loadcellwall, Neon

Use both single and double precision exe

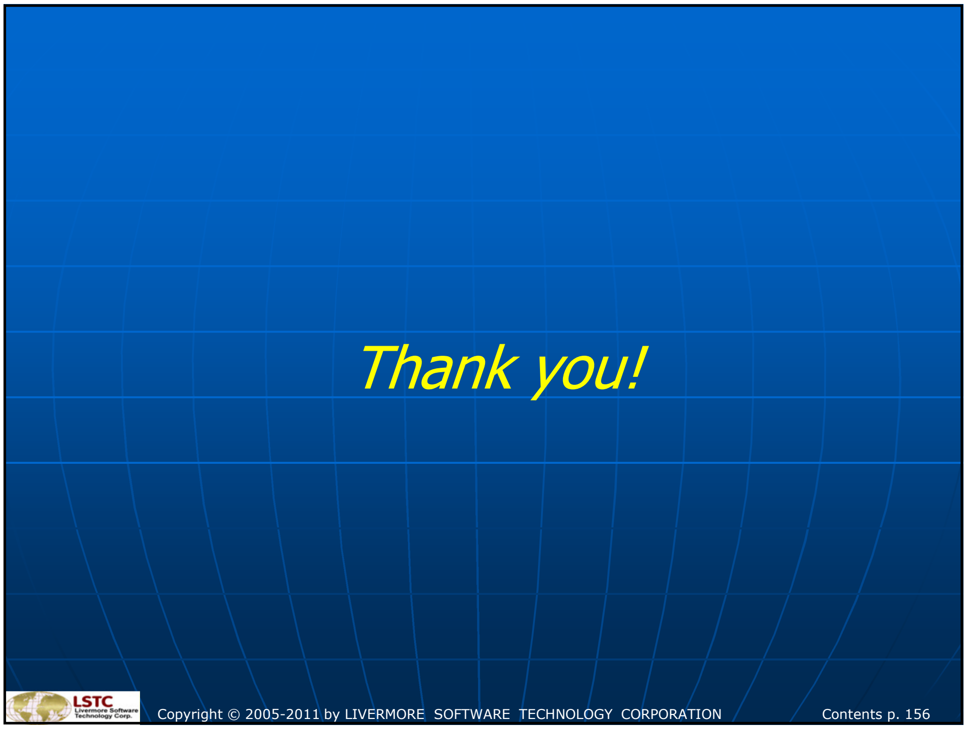


Neon Model Practice 2

E1: check system CPU info, NUMA order, memory, etc

E2: Decompose Neon into 16 partitions using double precision exe
use 4 processors single precision exe for analysis for 100 cycles
(use top to check CPU usage)

E3: use hybrid exe with 2 mpp and 2 threads each
(check the mpp971 usage from top command.)



Thank you!