Teradactyl: An Easy-to-Use Supercomputer

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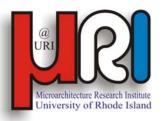
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- From: "*Pterodactyl*" winged fingers
 - Flying dinosaur
- *Teradactyl*: 'Flying' Supercomputer
 - Many 'fingers'
 - Each 'finger' (processing element) flies

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Work supported in present, past and future by:

• Laurette Bradley, my wife.

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- 1. The Problem
- 2. Our Approach to a Solution
- 3. Resource-Flow Execution
- 4. Basic LEVO Microarchitecture
- 5. Teradactyl Architecture
- 6. Summary

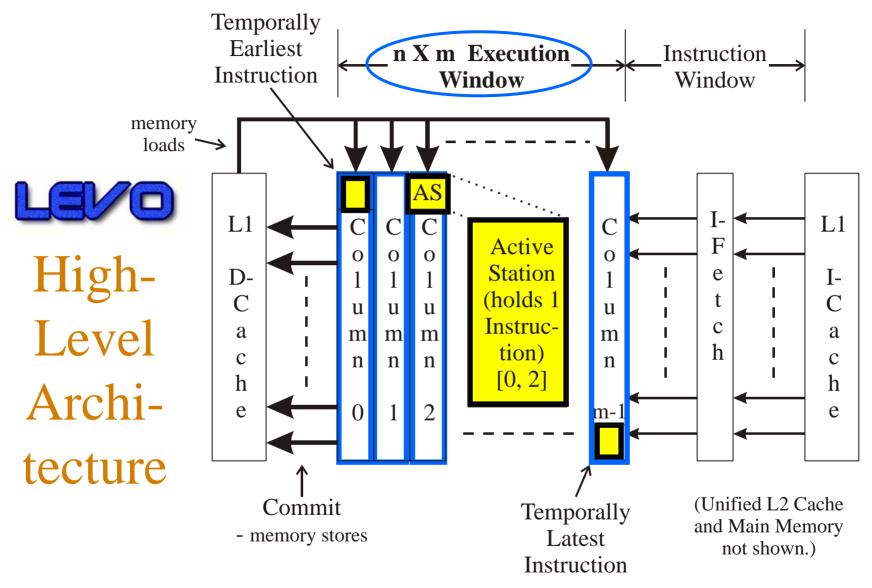


- PROGRAMMABILITY!!!
- Why? Here's why:
 - 1. Scientific programs by users take years to write
 - Even with libraries
 - "...the manual development and testing of a reasonably efficient parallel code for a computational model ... typically takes <u>months to years</u> for a computational chemist." (Our emphases.), Supercomputing 2002
 - Parallel programming, scheduling, etc. way too hard
 - Users' time is greatly misspent:
 chemists should be doing Chemistry, not coding

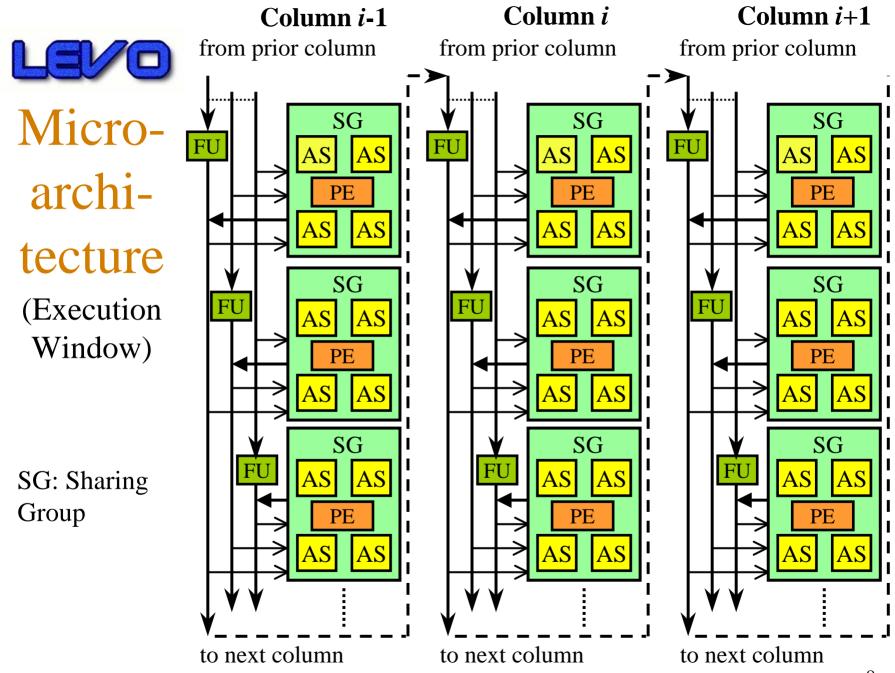


- 1. Use *hardware* to do the hard stuff
- Let the user use an *easy* programming model:
 → *Standard sequential (imperative) model*
- 3. WHY is parallel programming so hard?
 → don't know where the *data dependencies* are
- 4. Approach has always been to estimate them
- 5. Don't <u>estimate</u> them, <u>know</u> them
- 6. Use *resource-flow* execution:
 - Instructions flow to PE's, are executed <u>regardless</u> of dependencies
 - Then clean up: <u>enforce</u> dependencies when they're <u>known</u>

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Processing Elements (PEs) are distributed among AS's.







- Small integers = position in E-window
- <u>Enforce</u> and <u>minimize</u> dependencies
- Provide operand linking (sink-to-source)
- Basic problem:
 - R3 must wind up with *closest previous value* of R4 (2)
 - Must be independent of execution order of instructions

nominal time order
Instructions:
$$R4 = 1$$
 $R4 = 2$ $R3 = R4$
 $I1$ $I5$ $I9$
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• LSTT: Last Snarfed Time-Tag

nominal time orderInstructions:R4 = 1
I 1R4 = 2
I 5R3 = R4
I 9

 $\underline{Time - 1}$:I 1 brdcsts.R4 address matches,TT(I 1) >= LSTT(I 9),I 1 info snarfed: R3=1

<u>Time - 2</u>:

I 5 brdcsts.

R4 address matches, $TT(I 5) \ge LSTT(I 9)$, I 5 info snarfed: R3=<u>2</u>

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• Recall: R3 \leftarrow *closest previous value* of R4 (2)

Time - 1:I 5 brdcsts.R4 address matches,
TT(I 5) >= LSTT(I 9),
I 5 info snarfed: R3=2

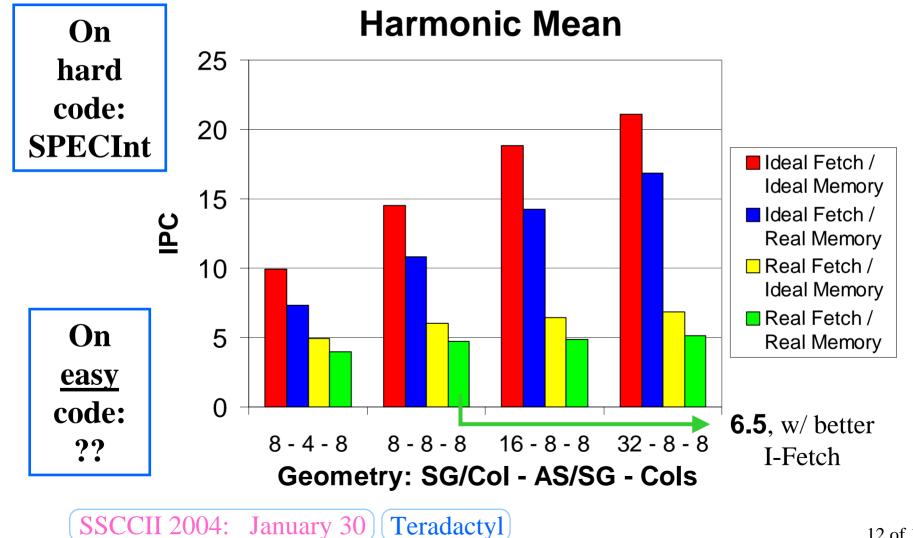
<u>*Time* – 2</u>: I 1 brdcsts.

R4 address matches, TT(I 1) < LSTT(I 9), I 1 info **not** snarfed.

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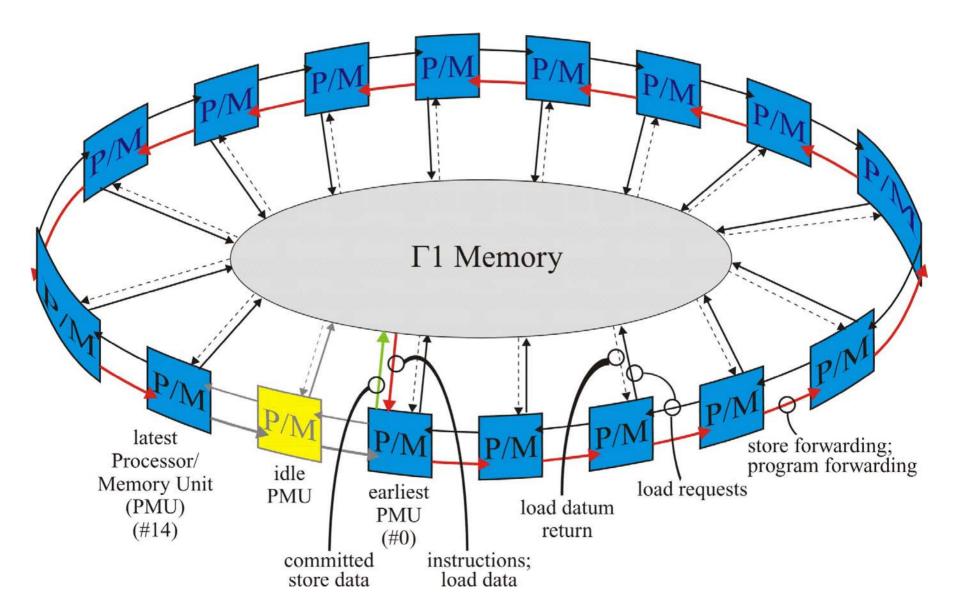






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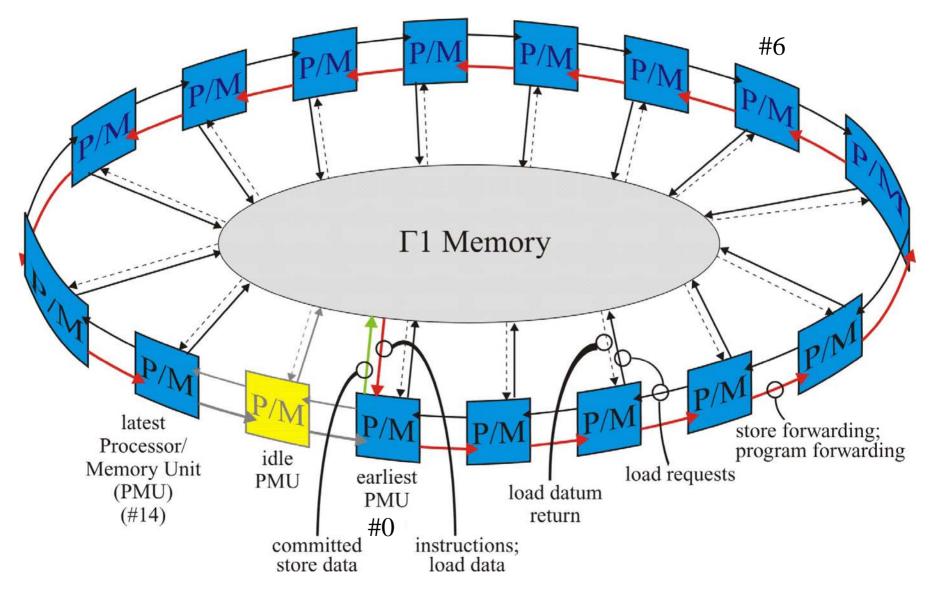






 $(A \times B) + C \qquad \frac{Before:}{After:} + (\#6) \text{ wait for 'x' (\#0); manual} \\ \underline{After:} + (\#6) \parallel (x' (\#0)); \& \text{ auto-schedule} \\ \end{array}$

Example





- Scalable to thousands of processors
- Uses modified LEVO model:
 - Processor Memory Units like Levo columns
 - Augment time-tags with PMU # as a prefix
 - Whole Teradactyl like Levo Execution Window
- Once data computed, is sent around ring, to update dependent operands:
 → close to best performance possible
 (→ with speculation, maybe better)



- First: Talking about <u>SUSTAINED</u> performance
- Now, assume:
 - PMU (Levo) up to 10 IPC within several years
 - Chip clock frequency up to 5 GHz \rightarrow 50 Gops
 - Then for TeraOp: ~25 PMU's (some inefficiency)
 And for PetaOp: 25,000 PMU's
- Power, etc.: ~ same as other supercomputers
- Other supercomputers: ~< 1 TeraOp, sustained
 With <u>100's</u> or <u>1000's</u> of processors





- Problem: Programmability
- Solution: <u>Teradactyl</u>:
 - Based on resource-flow execution
 - Data dependencies known exactly, at run-time
 - Data speculation also used to improve performance
 - Scalable
 - "Easy" to program (well, as easy as it can be :-)
- The future: <u>Petadactyl</u>

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