Introduction to Parallel Computing (CMSC416 / CMSC616)



Performance Modeling, Analysis, and Tools

Abhinav Bhatele, Department of Computer Science



Annoucements

- Assignment I is now posted online
 - Due on: Oct 2, 2023 11:59 pm
- More resources:
 - https://www.cs.umd.edu/~mmarsh/books/cmdline/cmdline.html
 - https://www.cs.umd.edu/~mmarsh/books/tools/tools.html
- Late submission policy: submit up to one late day for a 20% penalty
 - For any other exceptions, you need to ask as early as possible, not on the day of the deadline
- Quiz I will be posted on Wed Sep 20, 2023 I I:59 pm
 - You'll have 24 hours to take it



Weak versus strong scaling

- Strong scaling: Fixed total problem size as we run on more processes
 - Sorting n numbers on 1 process, 2 processes, 4 processes, ...
- Weak scaling: Fixed problem size per process but increasing total problem size as we run on more processes
 - Sorting n numbers on I process
 - 2n numbers on 2 processes
 - 4n numbers on 4 processes



Amdahl's law

- Speedup is limited by the serial portion of the code
 - Often referred to as the serial "bottleneck"
- Lets say only a fraction f of the code can be parallelized on p processes

Speedup =
$$\frac{1}{(1-f)+f/p}$$

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Performance analysis

- Parallel performance of a program might not be what the developer expects
- How do we find performance bottlenecks?
- Performance analysis is the process of studying the performance of parallel code
- Identify why performance might be slow
 - Serial performance
 - Serial bottlenecks when running in parallel
 - Communication overheads

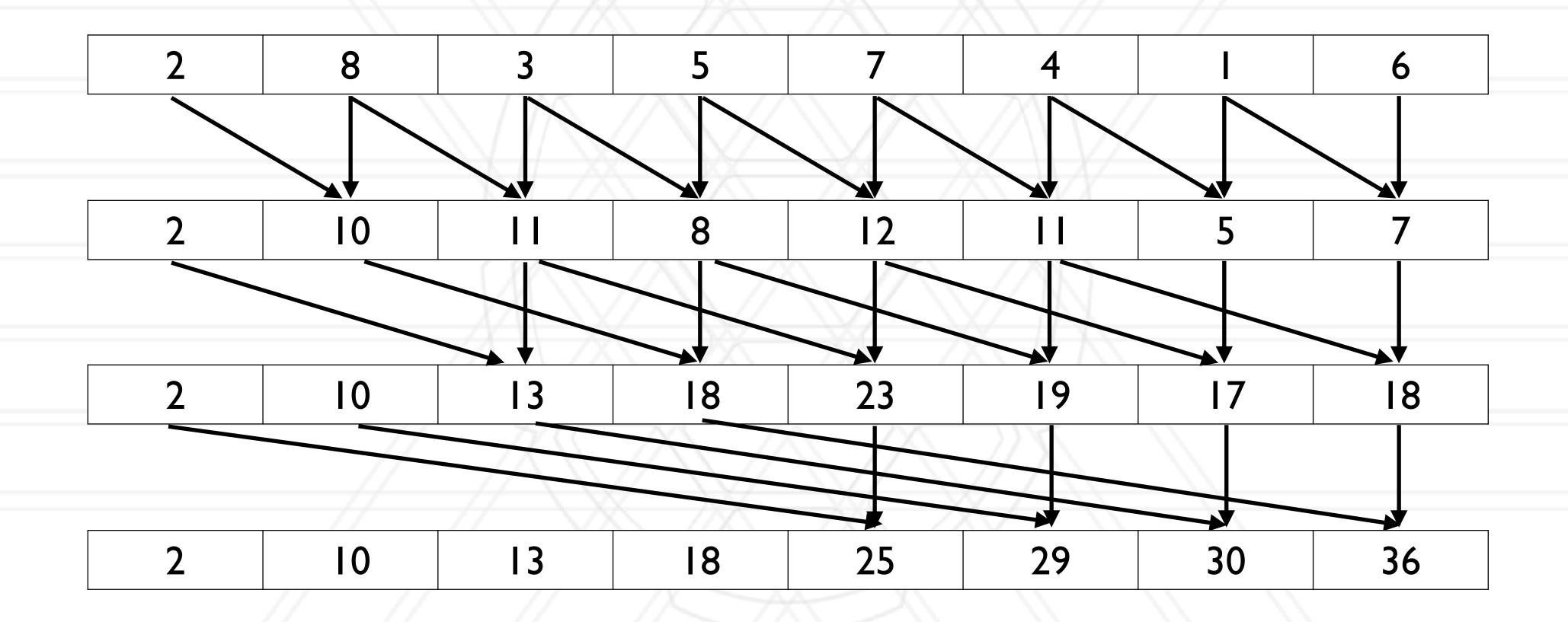


Performance analysis methods

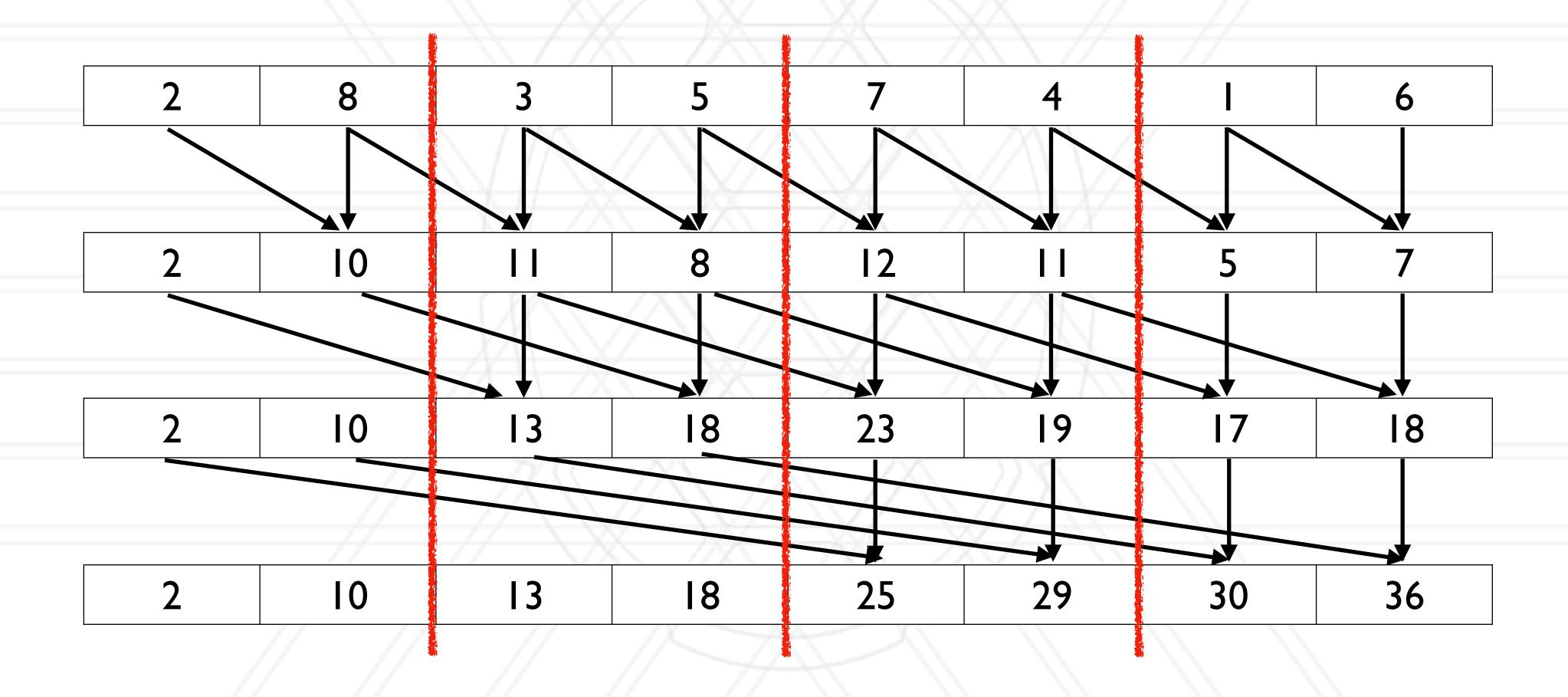
- Analytical techniques: use algebraic formulae
 - In terms of data size (n), number of processes (p)
- Time complexity analysis
- Scalability analysis (Isoefficiency)
- Model performance of various operations
 - Analytical models: LogP, alpha-beta model
- Empirical performance analysis using tools



Parallel prefix sum



Parallel prefix sum





- Assign n/p elements (block) to each process
- Perform prefix sum on these blocks on each process locally
 - Number of calculations:
- Then do parallel algorithm with partial prefix sums
 - Number of phases:
 - Total number of calculations:
 - Communication:



- Assign n/p elements (block) to each process
- Perform prefix sum on these blocks on each process locally
 - Number of calculations: $\frac{n}{p}$
- Then do parallel algorithm with partial prefix sums
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- Perform prefix sum on these blocks on each process locally
 - Number of calculations: $\frac{n}{p}$
- Then do parallel algorithm with partial prefix sums
 - Number of phases: log(p)
 - Total number of calculations: $log(p) \times \frac{n}{p}$
 - Communication:

Modeling communication: LogP model

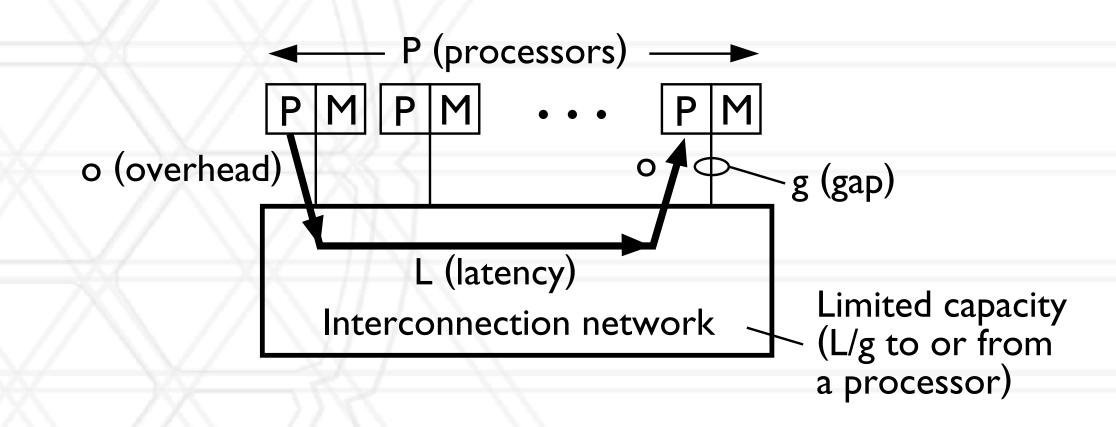
Model for communication on an interconnection network

L: latency or delay

o: overhead (processor busy in communication)

g: gap (between successive sends/recvs)

P: number of processors / processes





alpha + n * beta model

Another model for communication

$$T_{\text{comm}} = \alpha + n \times \beta$$

a: latency

n: size of message

I/β: bandwidth

Isoefficiency

- Relationship between problem size and number of processors to maintain a certain level of efficiency
- At what rate should we increase problem size with respect to number of processors to keep efficiency constant

Speedup and efficiency

• Speedup: Ratio of execution time on one process to that on p processes

Speedup =
$$\frac{t_1}{t_p}$$

Efficiency: Speedup per process

Efficiency =
$$\frac{t_1}{t_p \times p}$$

Efficiency in terms of overhead

 Total time spent in all processes = (useful) computation + overhead (extra computation + communication + idle time)

$$p \times t_p = t_1 + t_o$$

Efficiency =
$$\frac{t_1}{t_p \times p} = \frac{t_1}{t_1 + t_o} = \frac{1}{1 + \frac{t_o}{t_1}}$$

Isoefficiency function

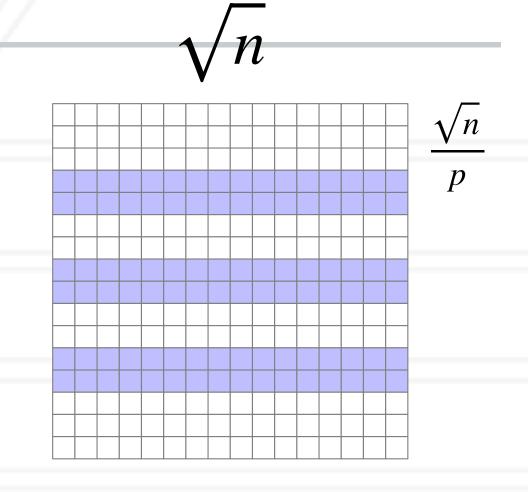
Efficiency =
$$\frac{1}{1 + \frac{t_o}{t_1}}$$

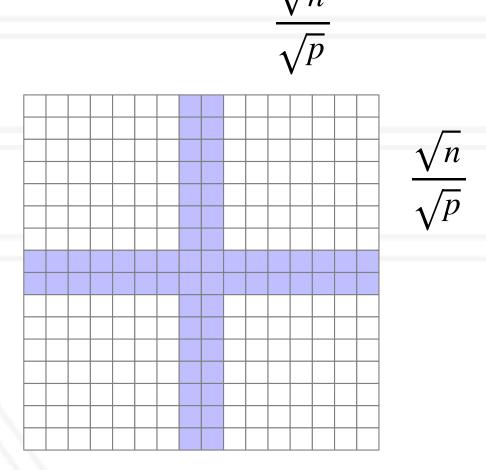
• Efficiency is constant if t_0 / t_1 is constant (K)

$$t_o = K \times t_1$$

- ID decomposition:
 - Computation:
 - Communication:

- 2D decomposition:
 - Computation:
 - Communication





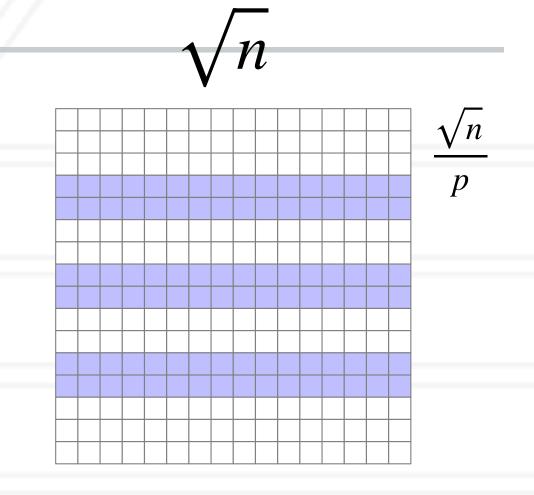
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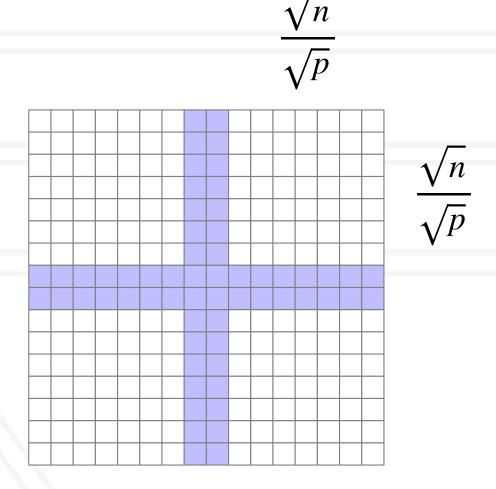
• Computation:
$$\sqrt{n} \times \frac{\sqrt{n}}{p} = \frac{n}{p}$$

Communication:



- Computation:
- Communication





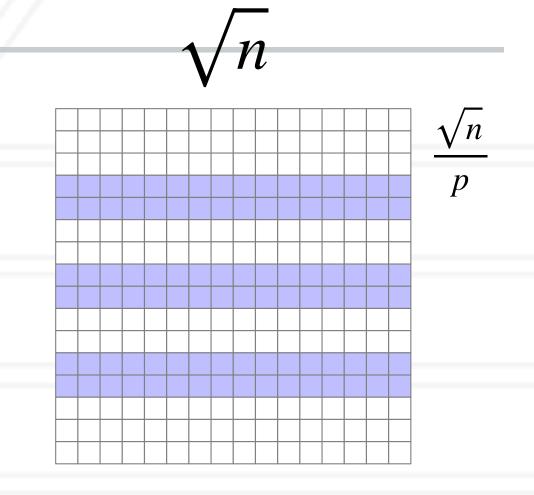
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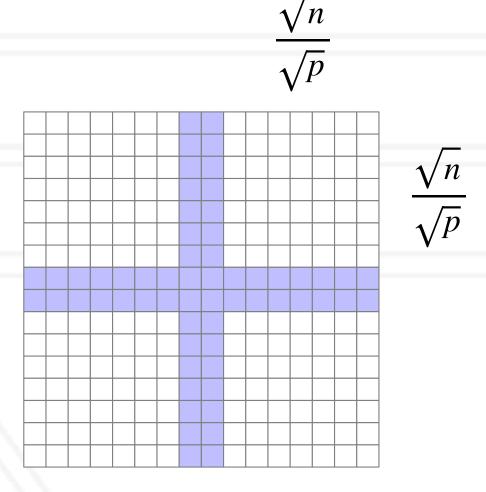
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• Communication: $2 \times \sqrt{n}$



- Computation:
- Communication

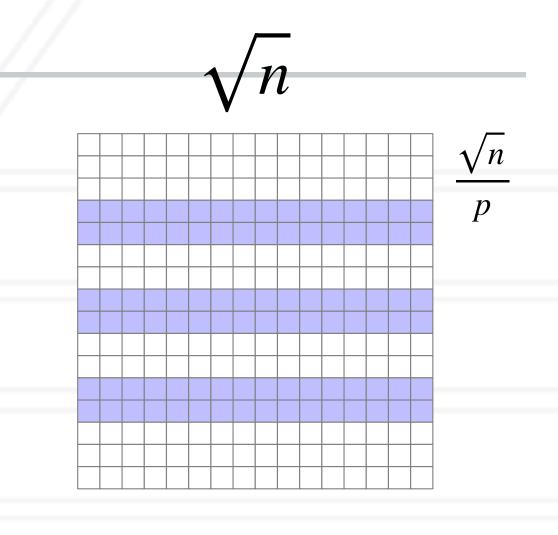




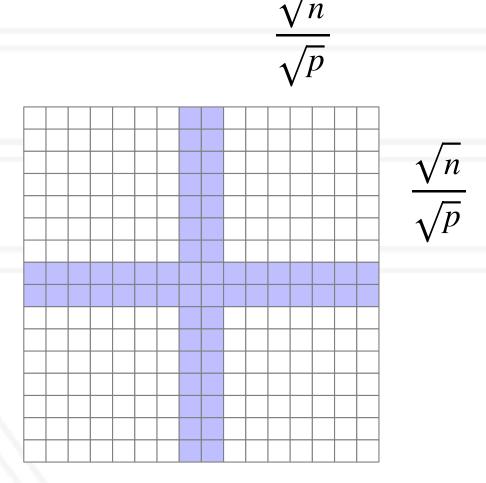
ID decomposition:

- Computation: $\sqrt{n} \times \frac{\sqrt{n}}{p} = \frac{n}{p}$
- Communication: $2 \times \sqrt{n}$

$$\frac{t_o}{t_1} = \frac{2 \times \sqrt{n}}{\frac{n}{p}} = \frac{2 \times p}{\sqrt{n}}$$



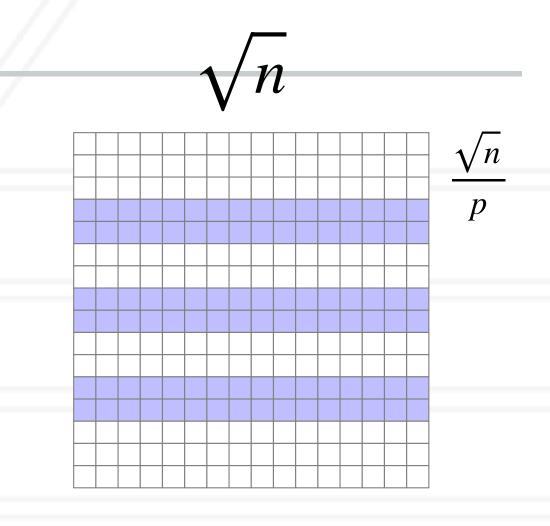
- 2D decomposition:
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ID decomposition:

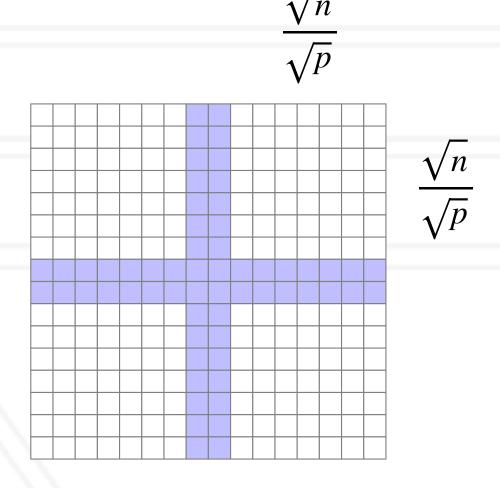
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• 2D decomposition:

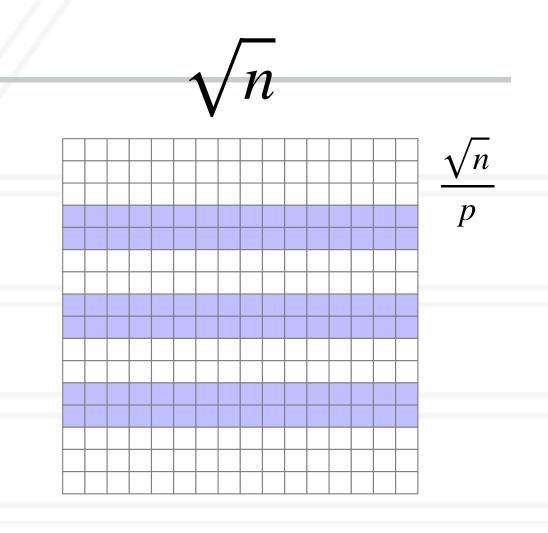
- Computation: $\frac{\sqrt{n}}{\sqrt{p}} \times \frac{\sqrt{n}}{\sqrt{p}} = \frac{n}{p}$
- Communication



ID decomposition:

- Computation: $\sqrt{n} \times \frac{\sqrt{n}}{p} = \frac{n}{p}$
- $2 \times \sqrt{n}$ Communication:

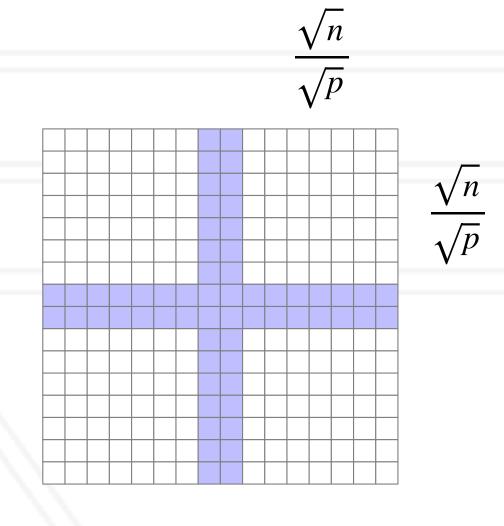
$$\frac{t_o}{t_1} = \frac{2 \times \sqrt{n}}{\frac{n}{p}} = \frac{2 \times p}{\sqrt{n}}$$



• 2D decomposition:

• Computation:
$$\frac{\sqrt{n}}{\sqrt{p}} \times \frac{\sqrt{n}}{\sqrt{p}} = \frac{n}{p}$$
• Communication
$$4 \times \frac{\sqrt{n}}{\sqrt{p}}$$

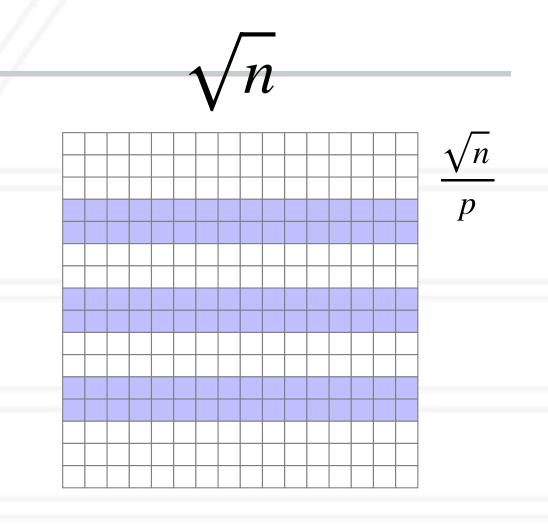
$$4 \times \frac{\sqrt{n}}{\sqrt{p}}$$



ID decomposition:

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- $2 \times \sqrt{n}$ Communication:

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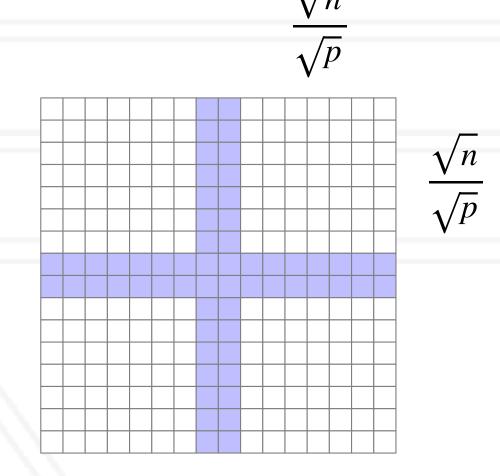
• 2D decomposition:

• Computation:
$$\frac{\sqrt{n}}{\sqrt{p}} \times \frac{\sqrt{n}}{\sqrt{p}} = \frac{n}{p}$$

• Communication $4 \times \frac{\sqrt{n}}{\sqrt{p}}$

$$4 \times \frac{\sqrt{n}}{\sqrt{p}}$$

$$\frac{t_o}{t_1} = \frac{4 \times \frac{\sqrt{n}}{\sqrt{p}}}{\frac{n}{p}} = \frac{4 \times \sqrt{p}}{\sqrt{n}}$$



Empirical performance analysis

- Two parts to performance analysis
 - measurement
 - analysis/visualization
- Simplest tool: timers in the code and printf



Using timers

```
double start, end;
double phase1, phase2, phase3;
start = MPI_Wtime();
 ... phase1 code ...
end = MPI Wtime();
phase1 = end - start;
start = MPI Wtime();
 ... phase2 ...
end = MPI_Wtime();
phase2 = end - start;
start = MPI_Wtime();
 ... phase3 ...
end = MPI_Wtime();
phase3 = end - start;
```



Using timers

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phase2 = end - start;
start = MPI Wtime();
 ... phase3 ...
end = MPI Wtime();
phase3 = end - start;
```

Phase I took 2.45 s

Phase 2 took 11.79 s

Phase 3 took 4.37 s



Performance tools

- Tracing tools
 - Capture entire execution trace, typically via instrumentation
- Profiling tools
 - Provide aggregated information
 - Typically use statistical sampling
- Many tools can do both



Metrics recorded

- Counts of function invocations
- Time spent in code
- Number of bytes sent
- Hardware counters
- To fix performance problems we need to connect metrics to source code



Tracing tools

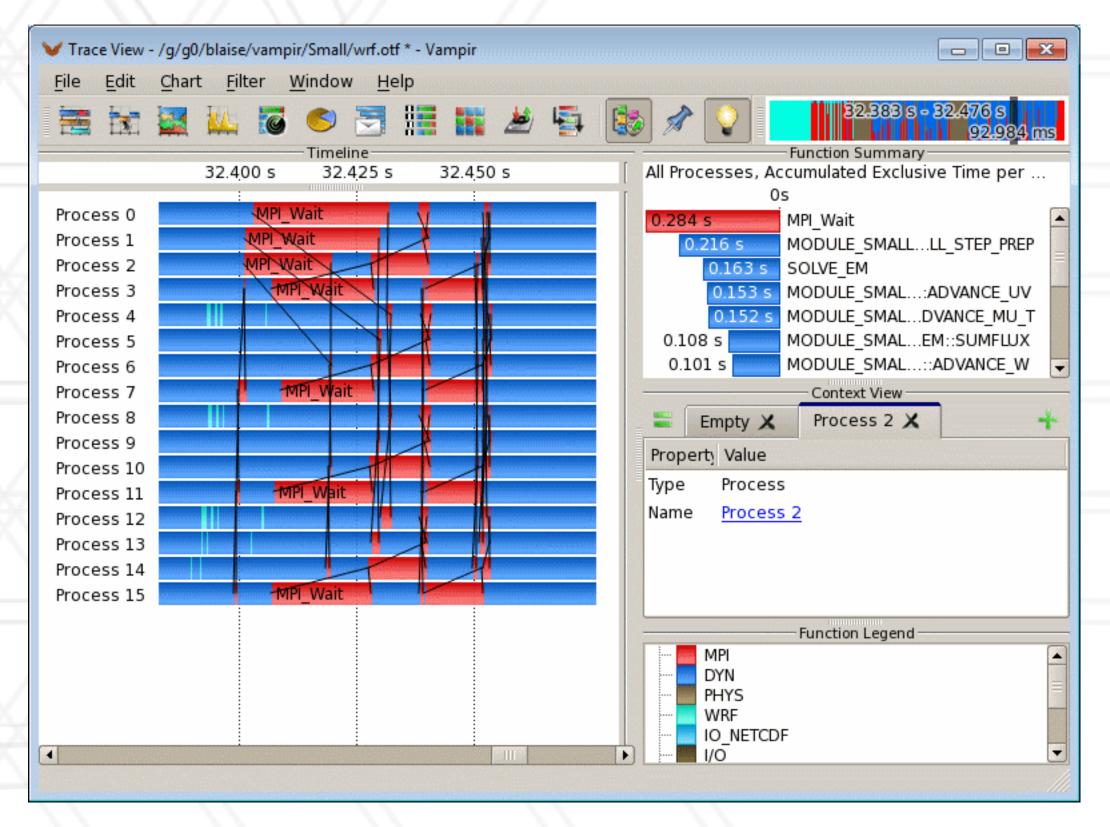
- Record all the events in the program with timestamps
- Events: function calls, MPI events, etc.

Vampir visualization: https://hpc.llnl.gov/software/development-environment-software/vampir-vampir-server



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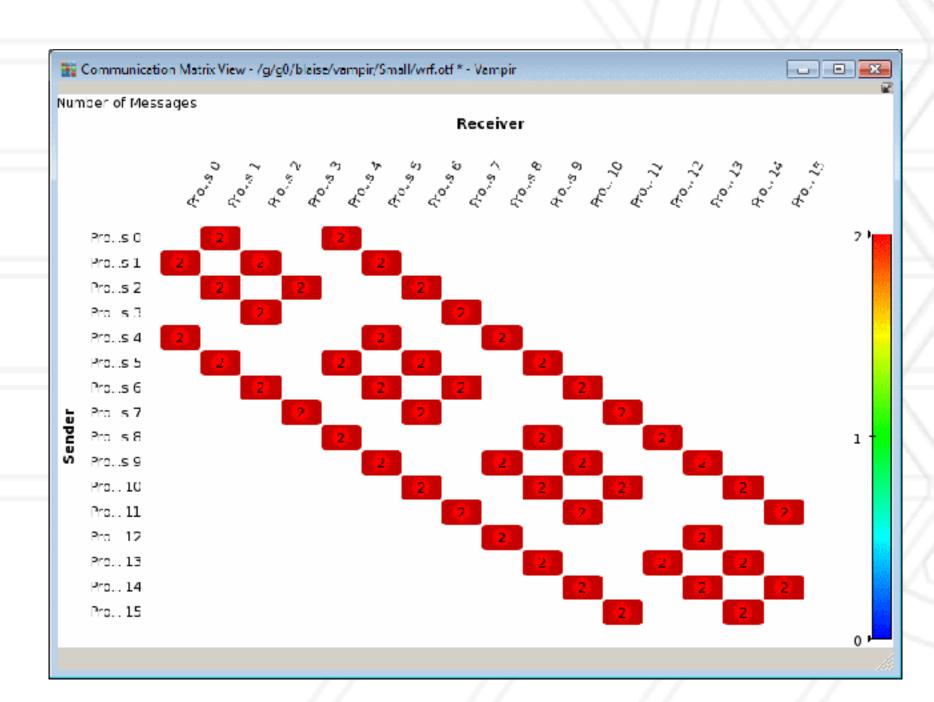


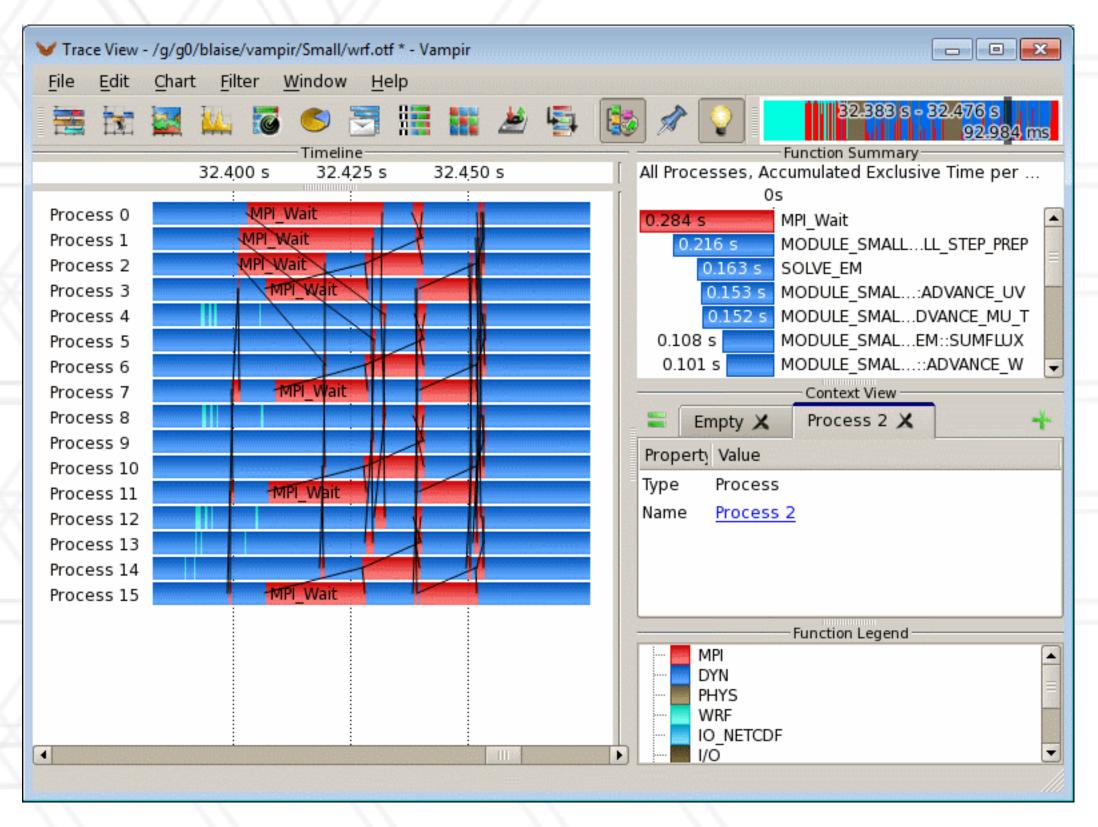
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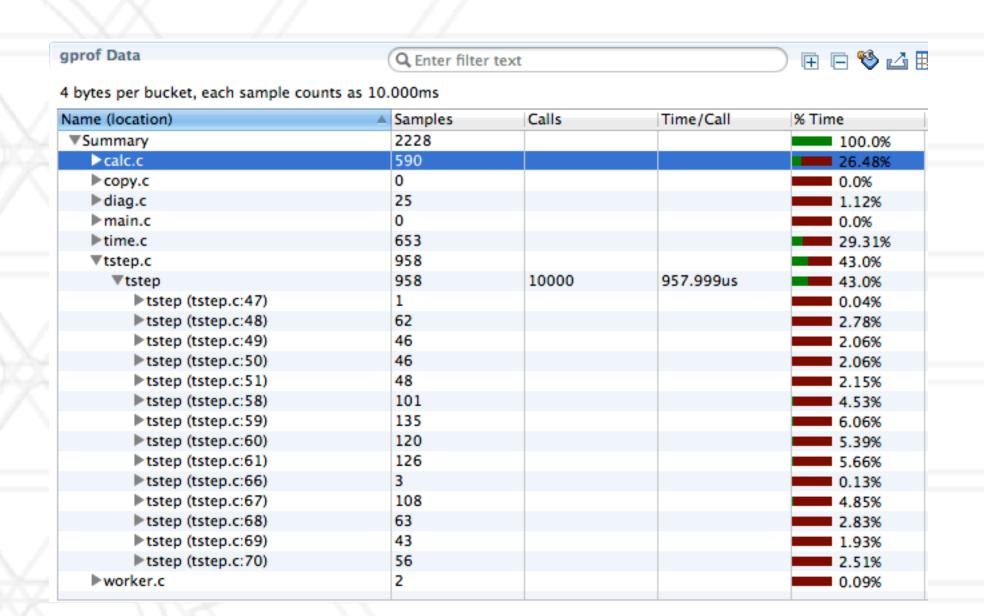
Examples of tracing tools

- VampirTrace
- Score-P
- TAU
- Projections
- HPCToolkit



Profiling tools

- Ignore the specific times at which events occurred
- Provide aggregate information about different parts of the code
- Examples:
 - Gprof, perf
 - mpiP
 - HPCToolkit, caliper
- Python tools: cprofile, pyinstrument, scalene

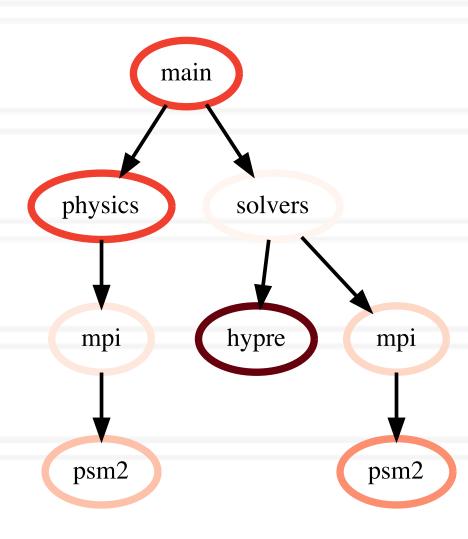


Gprof data in hpctView

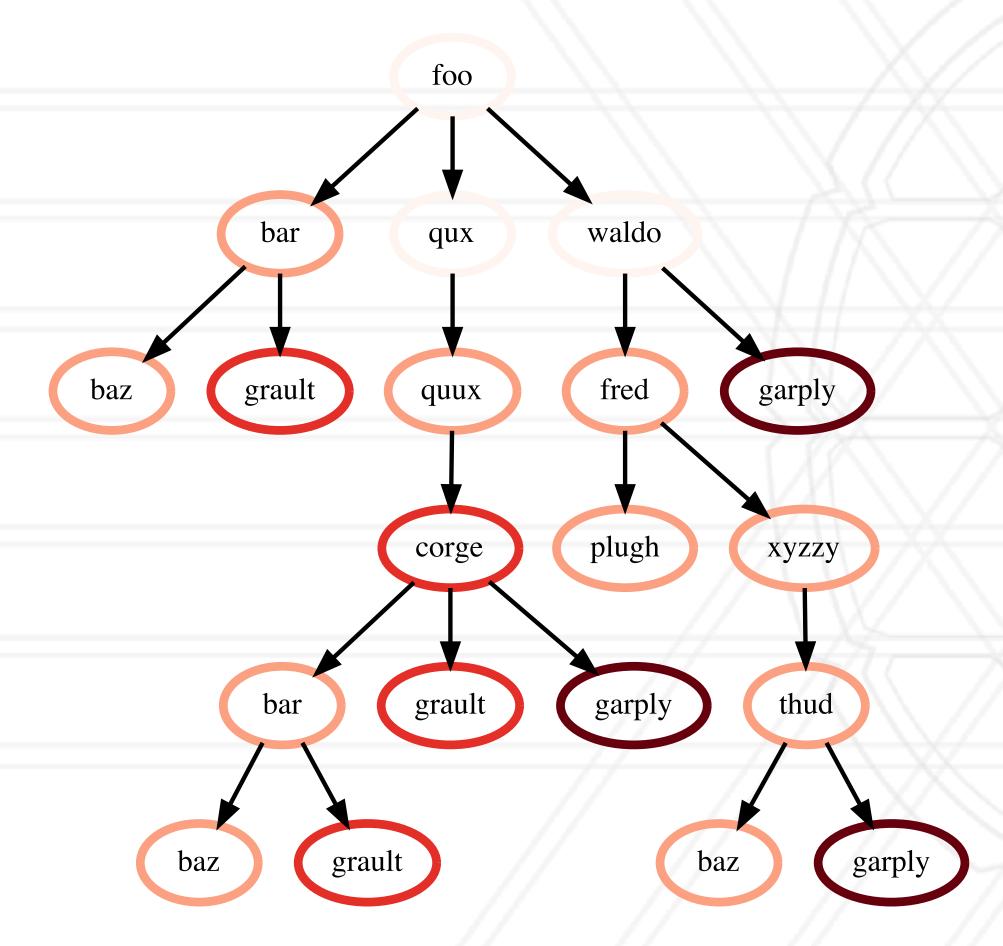


Calling contexts, trees, and graphs

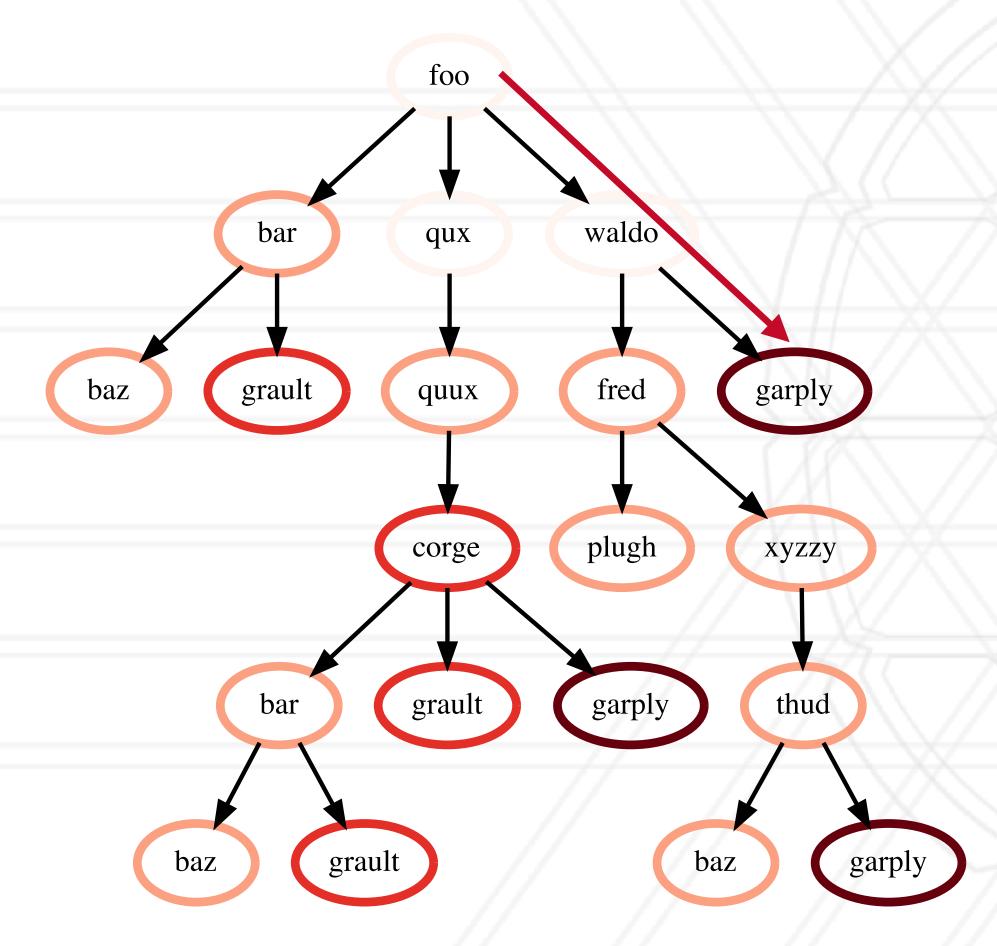
- Calling context or call path: Sequence of function invocations leading to the current sample
- Calling context tree (CCT): dynamic prefix tree of all call paths in an execution
- Call graph: merge nodes in a CCT with the same name into a single node but keep caller-callee relationships as arcs



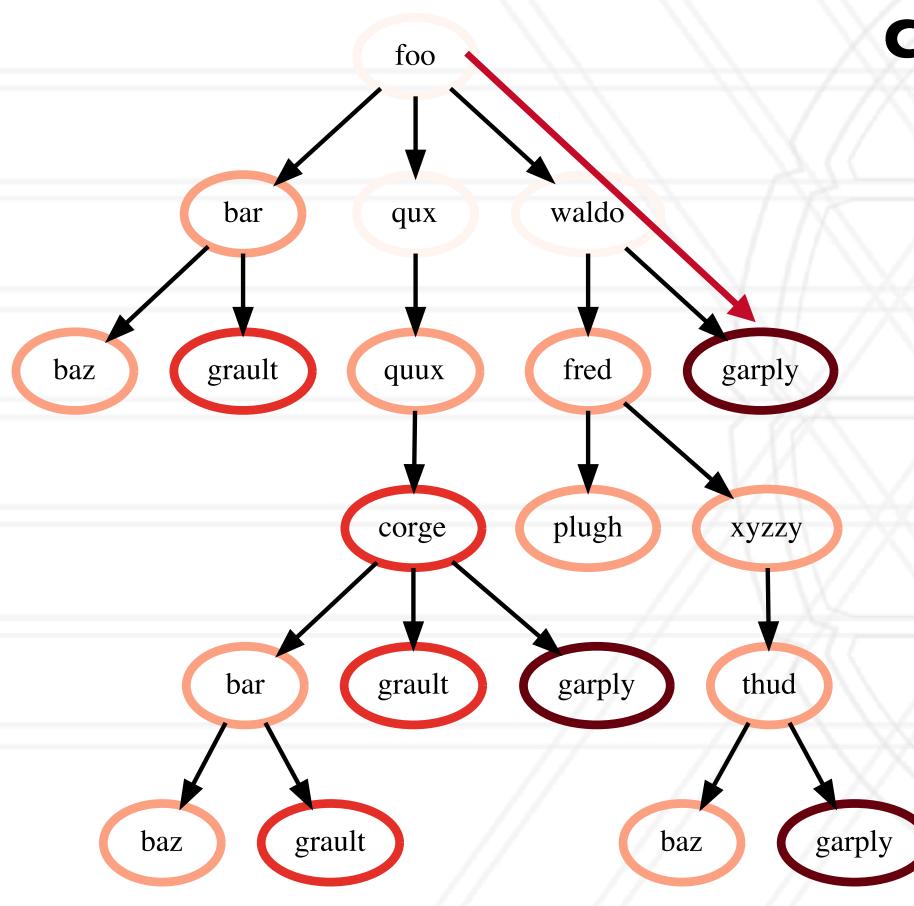








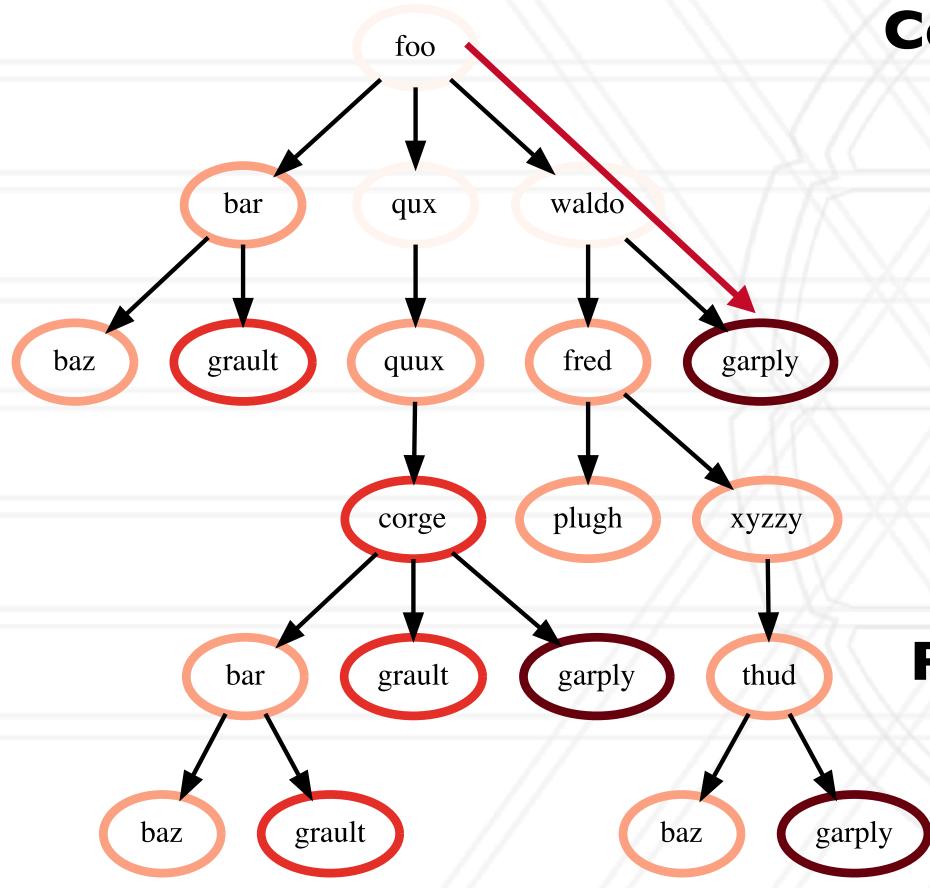




Contextual information

File
Line number
Function name
Callpath
Load module
Process ID
Thread ID





Contextual information

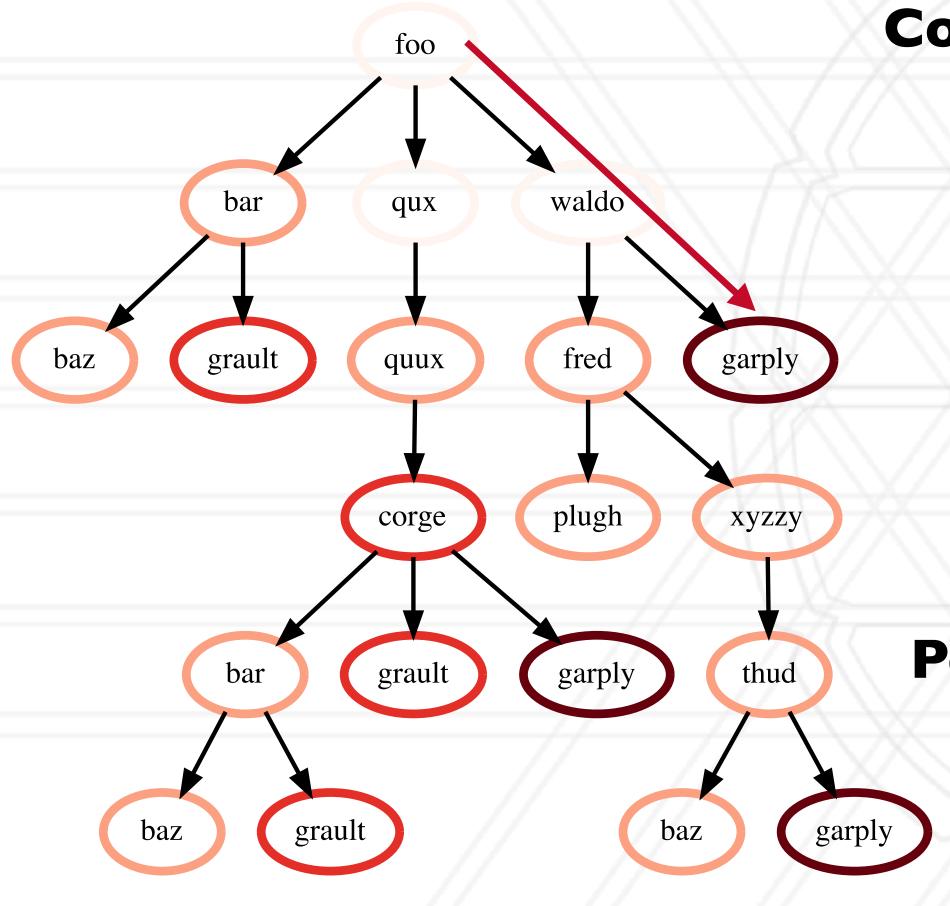
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Performance Metrics

Time
Flops
Cache misses





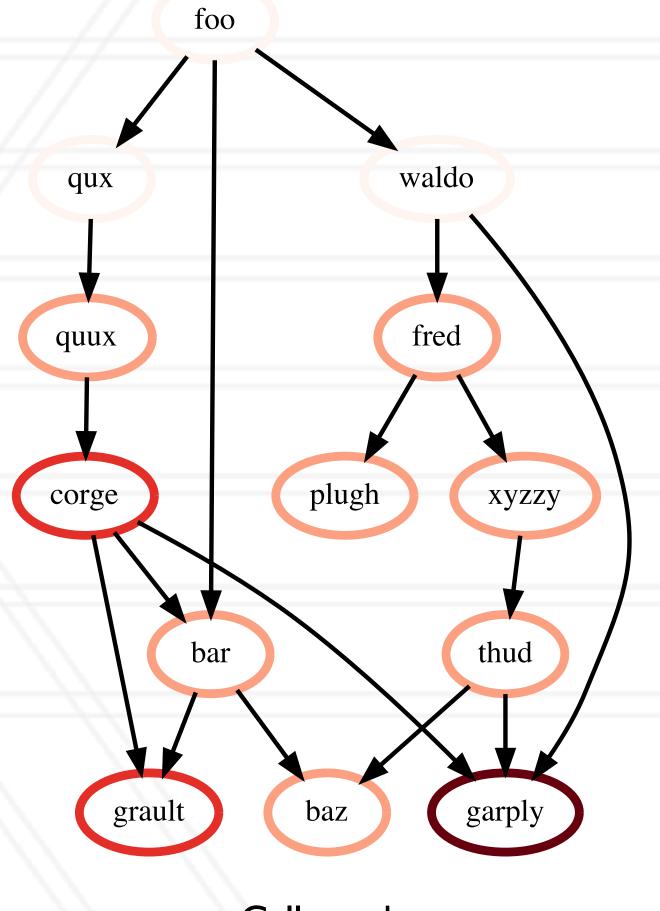
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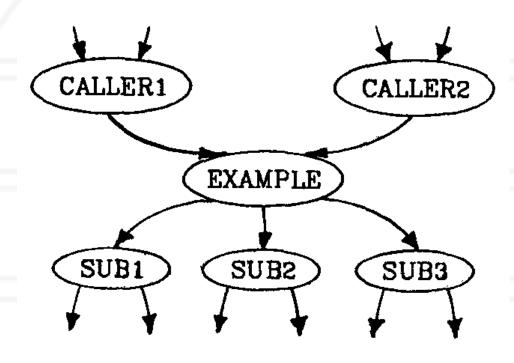


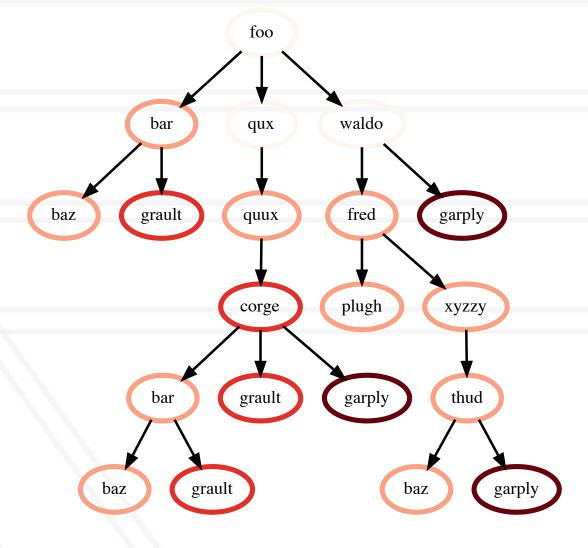


Call graph

Output of profiling tools

- Flat profile: Listing of all functions with counts and execution times
- Call graph profile
- Calling context tree







Hatchet

- Hatchet enables programmatic analysis of parallel profiles
- Leverages pandas which supports multi-dimensional tabular datasets
- Create a structured index to enable indexing pandas dataframes by nodes in a graph
- A set of operators to filter, prune and/or aggregate structured data

https://hatchet.readthedocs.io/en/latest/





 Pandas is an open-source Python library for data analysis

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- Dataframe: two-dimensional tabular data structure
 - Supports many operations borrowed from SQL databases

Columns

	node	name	time (inc)	time
0	{'name': 'main'}	main	200.0	10.0
1	{'name': 'physics'}	physics	60.0	40.0
2	{'name': 'mpi'}	mpi	20.0	5.0
3	{'name': 'psm2'}	psm2	15.0	30.0
4	{'name': 'solvers'}	solvers	100.0	10.0
5	{'name': 'hypre'}	hypre	65.0	30.0
6	{'name': 'mpi'}	mpi	35.0	20.0
7	{'name': 'psm2'}	psm2	25.0	60.0



Rows

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Rows

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- Dataframe: two-dimensional tabular data structure
 - Supports many operations borrowed from SQL databases
- Multilndex enables working with highdimensional data in a 2D data structure

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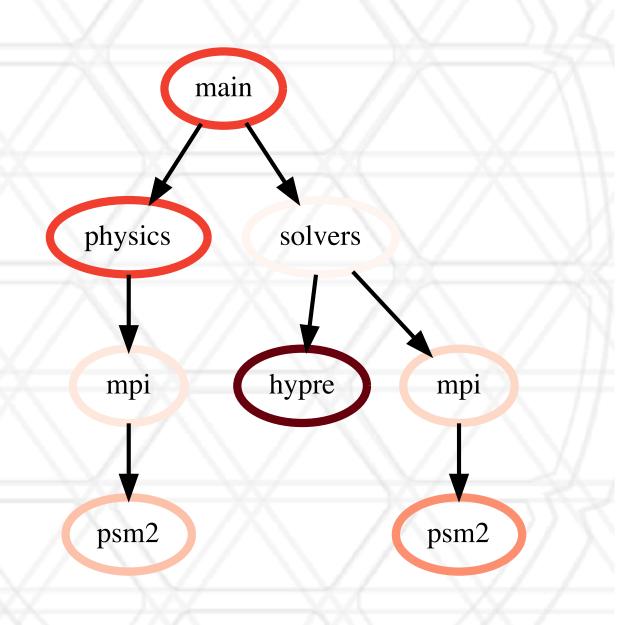
Central data structure: a GraphFrame

- Consists of a structured index graph object and a pandas dataframe
- Graph stores caller-callee relationships
- Dataframe stores all numerical and categorical data



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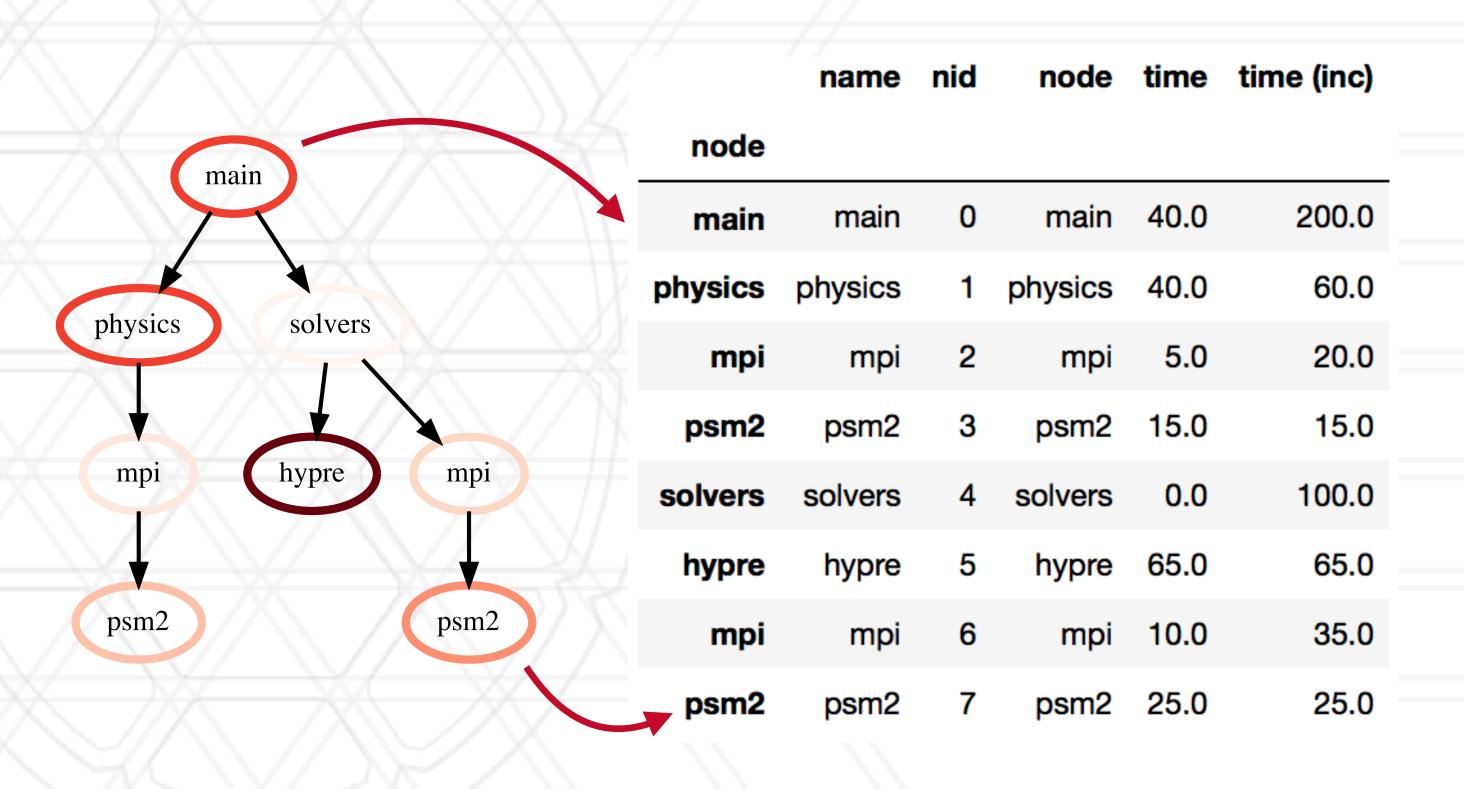
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Dataframe operation: filter

```
filtered_gf = gf.filter(lambda x: x['time'] > 10.0)
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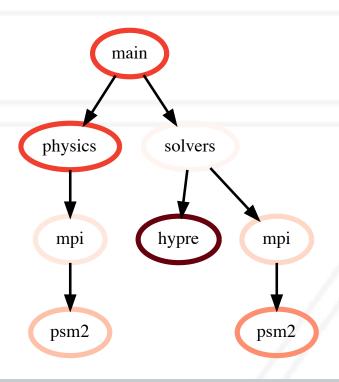
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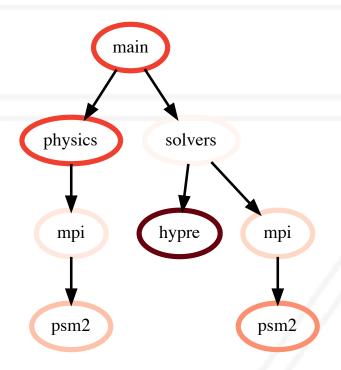


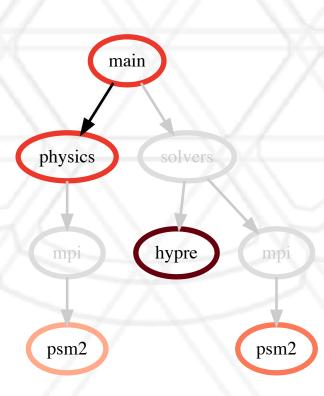
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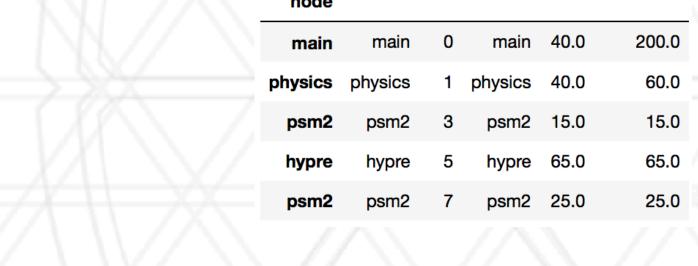


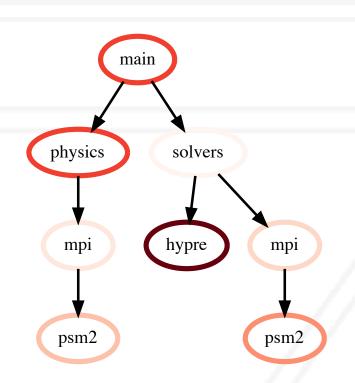
filter

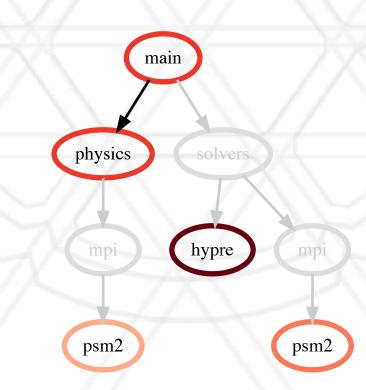
```
filtered_gf = gf.filter(lambda x: x['time'] > 10.0)
```

squashed_gf = filtered_gf.squash()

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node					
main	main	0	main	40.0	200.0
physics	physics	1	physics	40.0	60.0
mpi	mpi	2	mpi	5.0	20.0
psm2	psm2	3	psm2	15.0	15.0
solvers	solvers	4	solvers	0.0	100.0
hypre	hypre	5	hypre	65.0	65.0
mpi	mpi	6	mpi	10.0	35.0
psm2	psm2	7	psm2	25.0	25.0









filtered_gf = gf.filter(lambda x: x['time'] > 10.0)

squashed_gf = filtered_gf.squash()

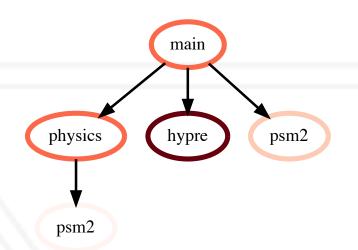
name	nid	node	time	time (inc)
main	0	main	40.0	200.0
physics	1	physics	40.0	60.0
mpi	2	mpi	5.0	20.0
psm2	3	psm2	15.0	15.0
solvers	4	solvers	0.0	100.0
hypre	5	hypre	65.0	65.0
mpi	6	mpi	10.0	35.0
psm2	7	psm2	25.0	25.0
	main physics mpi psm2 solvers hypre mpi	main 0 physics 1 mpi 2 psm2 3 solvers 4 hypre 5 mpi 6	main 0 main physics 1 physics mpi 2 mpi psm2 3 psm2 solvers 4 solvers hypre 5 hypre mpi 6 mpi	main 0 main 40.0 physics 1 physics 40.0 mpi 2 mpi 5.0 psm2 3 psm2 15.0 solvers 4 solvers 0.0 hypre 5 hypre 65.0 mpi 6 mpi 10.0

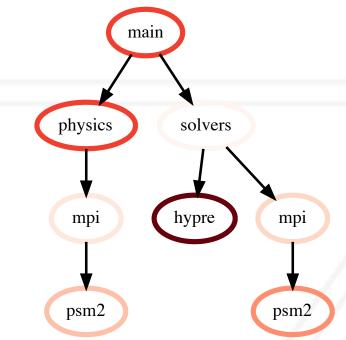


		name	nid	node	time	time (inc)
	node					
١	main	main	0	main	40.0	200.0
	physics	physics	1	physics	40.0	60.0
	psm2	psm2	3	psm2	15.0	15.0
	hypre	hypre	5	hypre	65.0	65.0
	psm2	psm2	7	psm2	25.0	25.0



	name	nıa	node	time	time (inc)
node					
main	main	0	main	40.0	200.0
physics	physics	1	physics	40.0	60.0
psm2	psm2	3	psm2	15.0	15.0
hypre	hypre	5	hypre	65.0	65.0
psm2	psm2	7	psm2	25.0	25.0





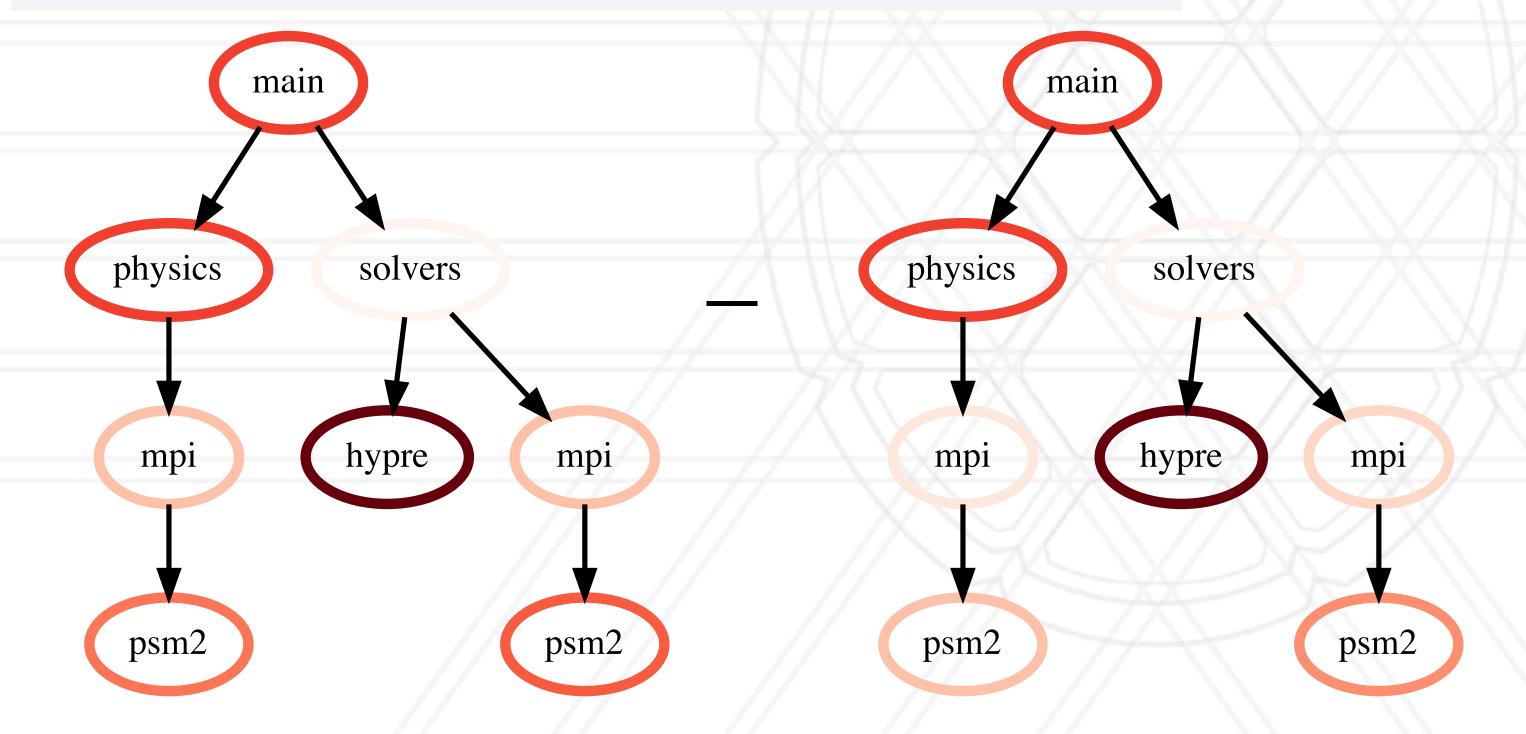


psm2

```
gf1 = ht.GraphFrame.from_literal( ... )
gf2 = ht.GraphFrame.from_literal( ... )
gf2 -= gf1
```

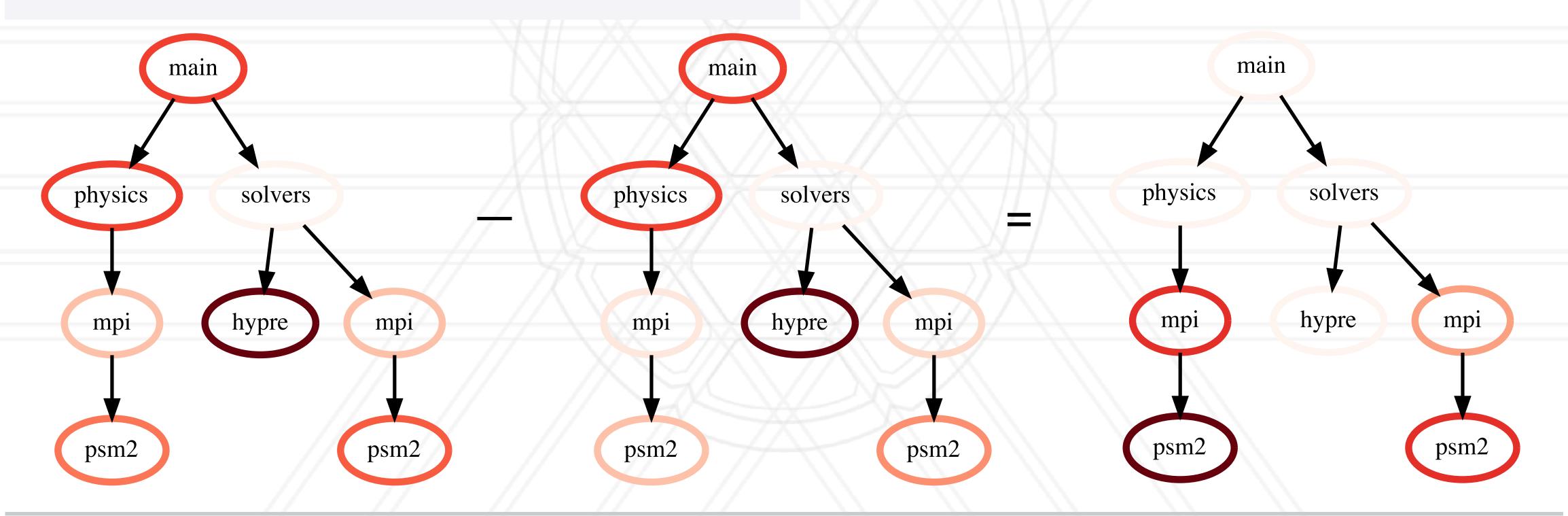


```
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```



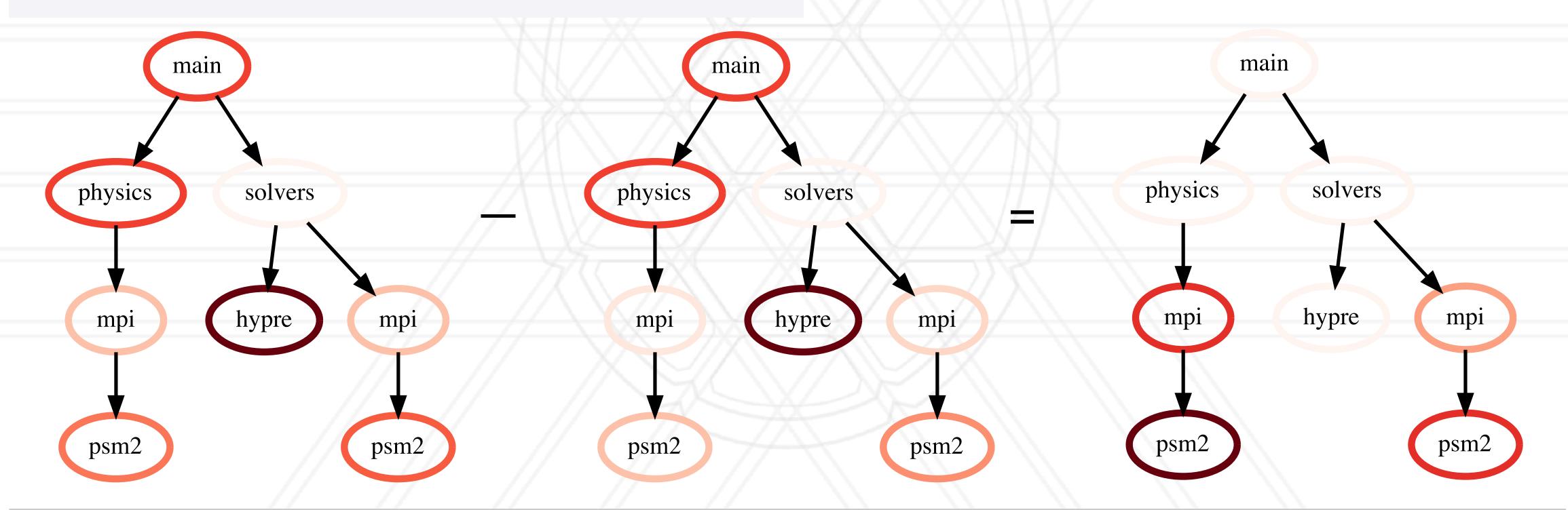


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```

https://hatchet.readthedocs.io



Visualizing small graphs

```
print(gf.tree(color=True))
```

```
0.000 foo
⊢ 5.000 bar
 └ 10.000 grault
├ 0.000 qux
  └ 5.000 quux
    └ 10.000 corge
      ⊢ 5.000 bar
       └ 10.000 grault
      └ 15.000 garply
└ 0.000 waldo
  ⊢ 5.000 fred
    └ 5.000 xyzzy
      └ 5.000 thud
        ⊢ 5.000 baz
        └ 15.000 garply
```

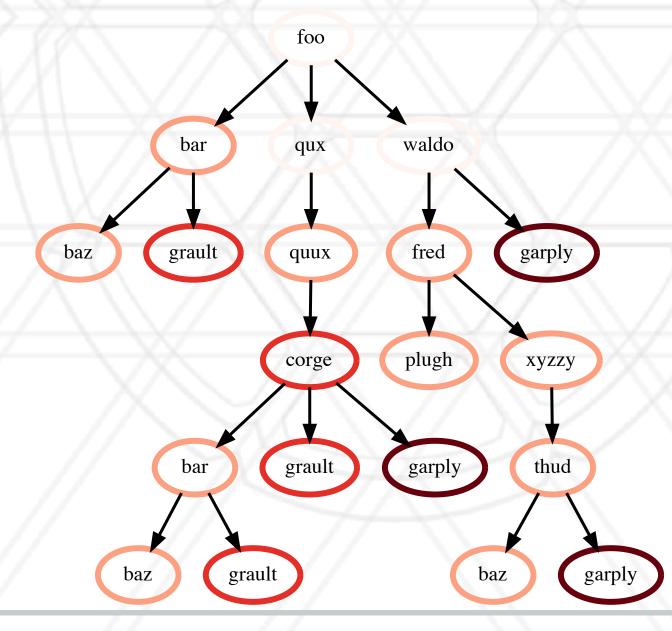


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0.000 foo
⊢ 5.000 bar
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  └ 10.000 grault
├ 0.000 qux
  └ 5.000 quux
     └ 10.000 corge
       ⊢ 5.000 bar
        ⊢ 5.000 baz
         └ 10.000 grault
       └ 15.000 garply
└ 0.000 waldo
  ⊢ 5.000 fred
    └ 5.000 xyzzy
       └ 5.000 thud
         ⊢ 5.000 baz
         └ 15.000 garply
```

```
with open("test.dot", "w") as dot_file:
    dot_file.write(gf.to_dot())
```





Visualizing small graphs

```
print(gf.tree(color=True))
```

```
0.000 foo
                                                                          with open("test.txt", "w") as folded_stack:
⊢ 5.000 bar
                            with open("test.dot", "w") as dot_file:
  ⊢ 5.000 baz
                                                                              folded_stack.write(gf.to_flamegraph())
   └ 10.000 grault
                                dot_file.write(gf.to_dot())
├ 0.000 qux
   └ 5.000 quux
     └ 10.000 corge
        ⊢ 5.000 bar
                                                                                                     waldo
          ⊢ 5.000 baz
                                                                                                         fred
                                                                                    quux
           └ 10.000 grault
                                                                                          corge
                                                                                                             XYZZY
        thud
        └ 15.000 garply
                                                         fred
                                                   quux
└ 0.000 waldo
   ⊢ 5.000 fred
                                                         plugh
      \vdash 5.000 plugh
                                                                                         Flamegraph
      └ 5.000 xyzzy
        └ 5.000 thud
           ⊢ 5.000 baz
           └ 15.000 garply
                                                grault
                                           baz
                                                             baz
```



Example 1: Generating a flat profile

```
gf = ht.GraphFrame.from_hpctoolkit('kripke')
gf.drop_index_levels()

grouped = gf.dataframe.groupby('name').sum()
sorted_df = grouped.sort_values(by=['time'], ascending=False)
print(sorted_df)
```



Example 1: Generating a flat profile

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```

		nid	11	time	time (inc)
E	name				
	<unknown file=""> [kripke]:0</unknown>	17234	1.8252826	80+ s	1.825282e+08
	Kernel_3d_DGZ::scattering	60	7.66993	e+07	7.896253e+07
	Kernel_3d_DGZ::LTimes	30	5.0104396	e+07	5.240528e+07
	Kernel_3d_DGZ::LPlusTimes	115	4.9477076	e+07	5.104498e+07
	Kernel_3d_DGZ::sweep	981	5.0188626	e+06	5.018862e+06
	memset.S:99	3773	3.1689826	e+06	3.168982e+06
	memset.S:101	3970	2.1208956	+06	2.120895e+06
	Grid_Data::particleEdit	1201	1.1312656	+06	1.249157e+06
	<unknown file=""> [libpsm2.so.2.1]:0</unknown>	324763	9.7334156	+05	9.733415e+05
	memset.S:98	3767	6.19777	e+05	6.197776e+05



Example 2: Comparing two executions

```
gf1 = ht.GraphFrame.from_caliper('lulesh-1core.json')
gf2 = ht.GraphFrame.from_caliper('lulesh-27cores. json')

gf2.drop_index_levels()
gf3 = gf2 - gf1

sorted_df = gf3.dataframe.sort_values(by=['time'], ascending=False)
print(sorted_df)
```



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```



Example 2: Comparing two executions

TimeIncrement	TimeIncrement	25.0	8.505048e+06	8.505048e+06
CalcQForElems	CalcQForElems	16.0	4.455672e+06	5.189453e+06
CalcHourglassControlForElems	CalcHourglassControlForElems	7.0	3.888798e+06	4.755817e+06
LagrangeNodal	LagrangeNodal	3.0	1.986046e+06	8.828475e+06
CalcForceForNodes	CalcForceForNodes	4.0	1.017857e+06	6.842429e+06



Example 3: Scaling study

```
datasets = glob.glob('lulesh*.json')
datasets.sort()
dataframes = []
for dataset in datasets:
   gf = ht.GraphFrame.from_caliper(dataset)
   gf.drop_index_levels()
    num_pes = re.match('(.*)-(\d+)(.*)', dataset).group(2)
    gf.dataframe['pes'] = num_pes
    filtered_gf = gf.filter(lambda x: x['time'] > 1e6)
    dataframes.append(filtered_gf.dataframe)
result = pd.concat(dataframes)
pivot_df = result.pivot(index='pes', columns='name', values='time')
pivot_df.loc[:,:].plot.bar(stacked=True, figsize=(10,7))
```

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datasets = glob.glob('lulesh*.json')
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```



Example 3: Scaling study

```
datasets = glob.glob('lulesh*.json')
    datasets.sort()
    dataframes = []
    for dataset in datasets:
                                                                                                       CalcEnergyForElems
         gf = ht.GraphFrame.from_caliper(datas
                                                                                                       CalcFBHourglassForceForElems
                                                                                                     CalcHourglassControlForElems

    CalcKinematicsForElems

         gf.drop_index_levels()
                                                                                                       CalcMonotonicOGradientsForElems
                                                                                                       CalcMonotonicQRegionForElems
                                                                                                     CalcPressureForElems
                                                                                                     EvalEOSForElems
         num_pes = re.match('(.*)-(\d+)(.*)',
                                                                                                     IntegrateStressForElems
                                                                                                     LagrangeNodal
         gf.dataframe['pes'] = num_pes
                                                                                                     MPI Allreduce
                                                                                                     MPI Waitall
         filtered_gf = gf.filter(lambda x: x['
         dataframes.append(filtered_gf.datafra
    result = pd.concat(dataframes)
pivot_df = result.pivot(index='pes', colu
    pivot_df.loc[:,:].plot.bar(stacked=True, figsize=(10,7))
```



Abhinav Bhatele

5218 Brendan Iribe Center (IRB) / College Park, MD 20742

phone: 301.405.4507 / e-mail: bhatele@cs.umd.edu