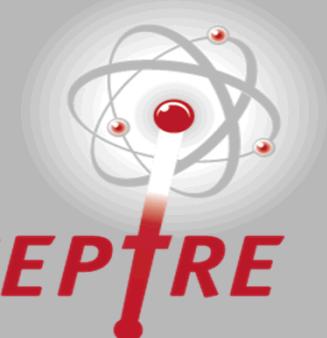


Parallel Deterministic Transport Sweeps of Structured and Unstructured Meshes with Overloaded Mesh Decompositions



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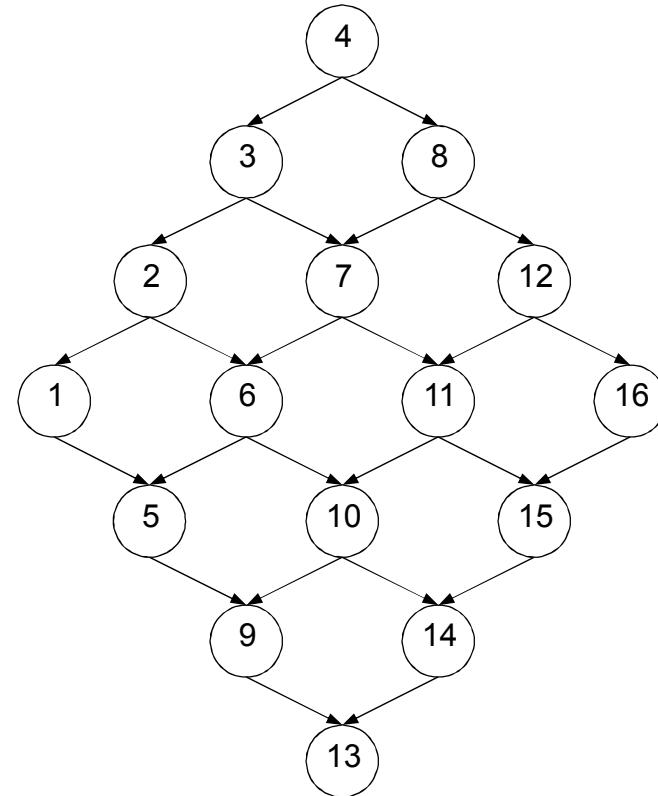
Outline

- Challenge for parallel scaling of sweeps
- General characterization of sweep algorithms
- Review of overloading technique for structured meshes
- Extension of structured overloading technique to unstructured meshes
- Results
- Conclusions

Challenge for parallel sweep scaling

$$\xrightarrow{\Omega}$$

1	2	3	4
5	6	7	8
9	10	11	12
13	14	15	16



Amount of concurrency < number of tasks

Challenge is to exploit existing concurrency

Characterization of sweep algorithms

Parallel sweep algorithms may be described by their approach to several issues

- Mesh partitioning
- Task aggregation
- Communication aggregation
- Task scheduling

Note that some approaches have used mesh partitioning to define task aggregation. Also, task and communication aggregation have typically been considered the same concept.

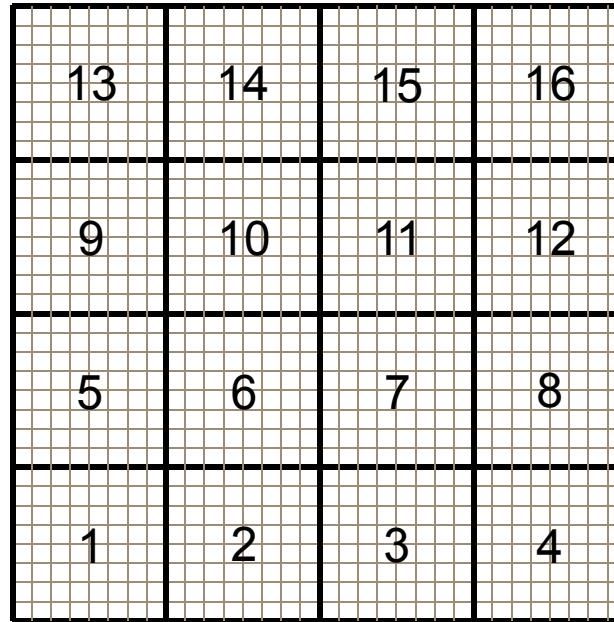
Structured vs. unstructured algorithms

Various parallel sweep techniques have been originally conceived and analyzed for structured meshes. Extending them to the unstructured case is often difficult.

We have recently extended the method of mesh overloading to the unstructured case.

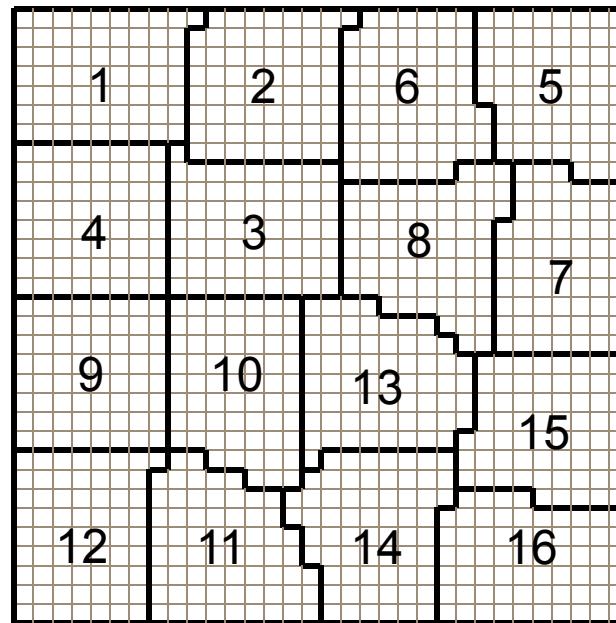
Note: “unstructured” algorithms may be applied to structured meshes, which is useful for comparison to structured algorithms.

Standard structured mesh partitioning



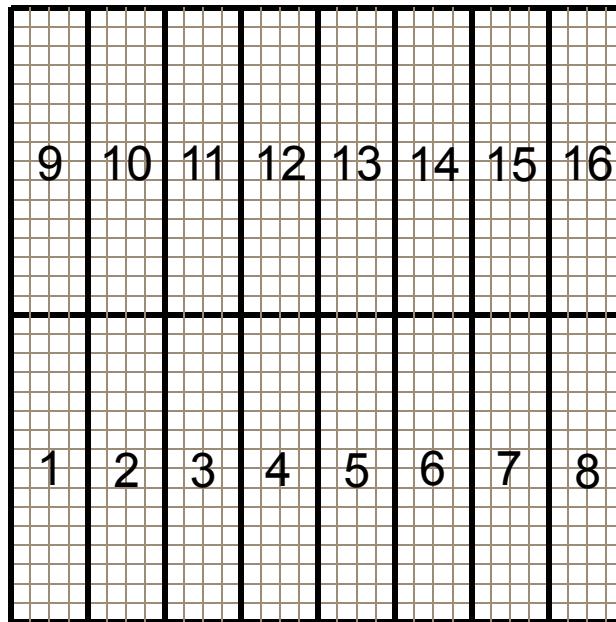
The standard approach to structured mesh partitioning subdivides the mesh into a $P_x \times P_y$ set of subdomains. It suffers from poor parallel sweep scaling.

Standard unstructured mesh partitioning



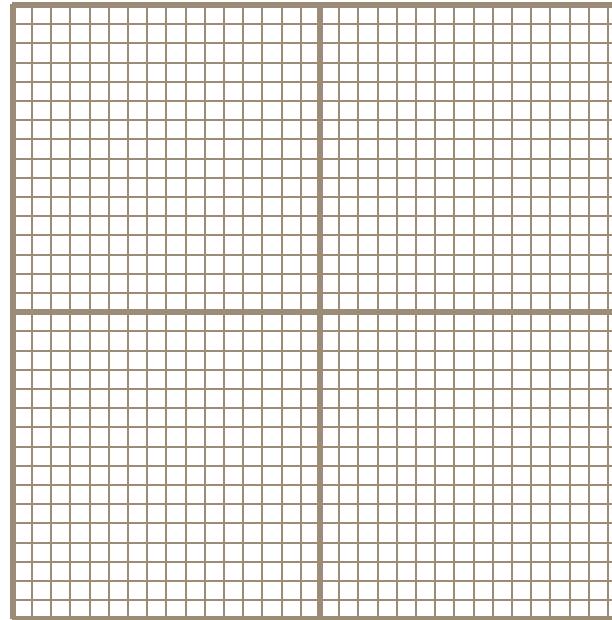
The standard approach to unstructured mesh partitioning subdivides the mesh into a set of P compact subdomains (e.g. with Chaco, Metis), which is a generalization of the structured approach. It also suffers from poor parallel sweep scaling.

Hybrid KBA partitioning



The hybrid KBA approach to structured mesh partitioning subdivides the mesh into a $P_x \times 2$ set of subdomains. It generally gives good parallel sweep scaling, but it is not clear how to extend it to unstructured meshes.

Structured mesh overloading



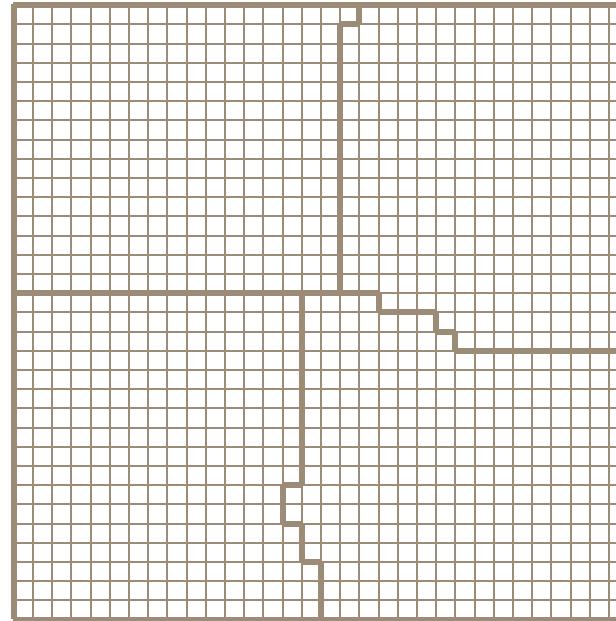
In the overloading approach for structured meshes, we first (conceptually) divide the mesh into a set of $\omega_x \times \omega_y$ “tiles”.

Structured mesh overloading, continued

13	14	15	16	13	14	15	16
9	10	11	12	9	10	11	12
5	6	7	8	5	6	7	8
1	2	3	4	1	2	3	4
13	14	15	16	13	14	15	16
9	10	11	12	9	10	11	12
5	6	7	8	5	6	7	8
1	2	3	4	1	2	3	4

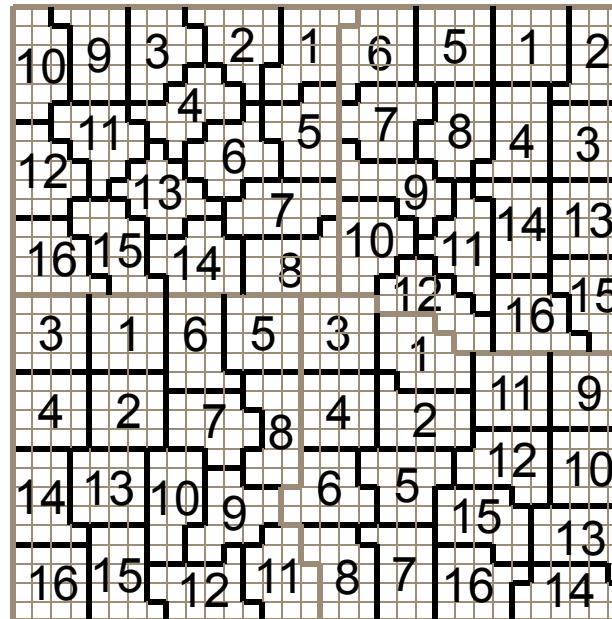
Next we divide each tile into a set of $P_x \times P_y$ subdomains. Each process gets a subdomain from each tile. This increases the sweep concurrency, but it also increases communication costs.

Unstructured mesh overloading



We extend the overloading approach to unstructured meshes by first dividing the mesh into a set of ω tiles by means of the usual mesh/graph partitioning algorithms.

Unstructured mesh overloading, continued



Next we divide each tile into a set of P subdomains. Each process gets a subdomain from each tile. This increases the sweep concurrency, but it also increases communication costs.

Advantages of unstructured mesh overloading

- Concepts from structured algorithms may be generalized to unstructured meshes
- May reuse existing algorithms created for standard unstructured mesh partitioning
- May balance computational and communication costs by means of the overloading parameter ω

Results

The Sceptre code was used to perform studies of the overloading approach for both structured and unstructured algorithms, in both cases applied to structured meshes. Sceptre itself is a general unstructured mesh code; it has no knowledge of structured properties. Partitioning occurs outside of the code as a preprocessing step.

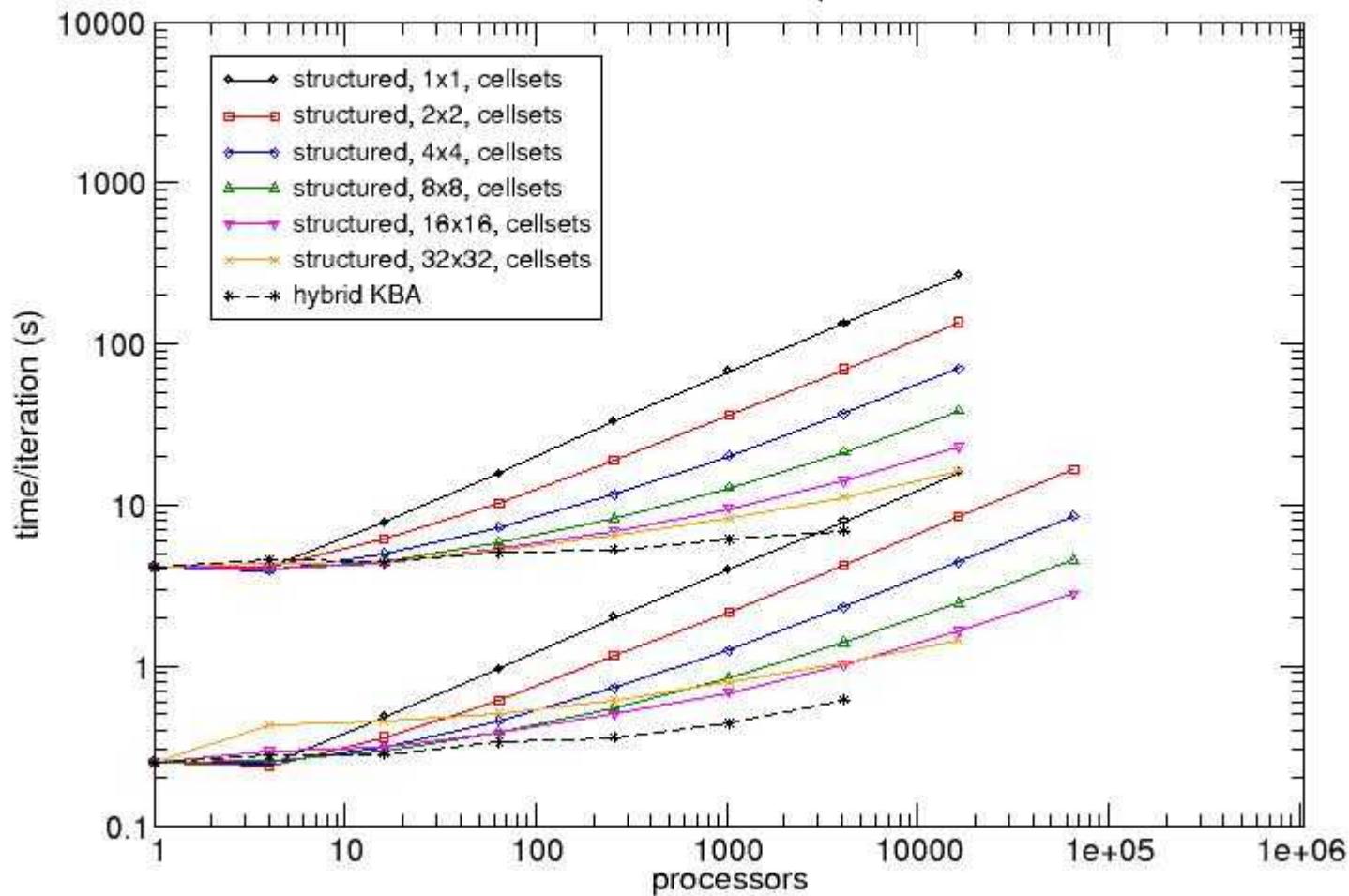
For the structured case we use cellsets that are identical to a full subdomain within a tile. For the unstructured case we use individual cells as a cellset. Communication occurs after the completion of each cellset.

We use the DFHDS scheduling heuristic in all cases.

All results were generated on Cielo, a massively parallel computer at Los Alamos National Laboratory.

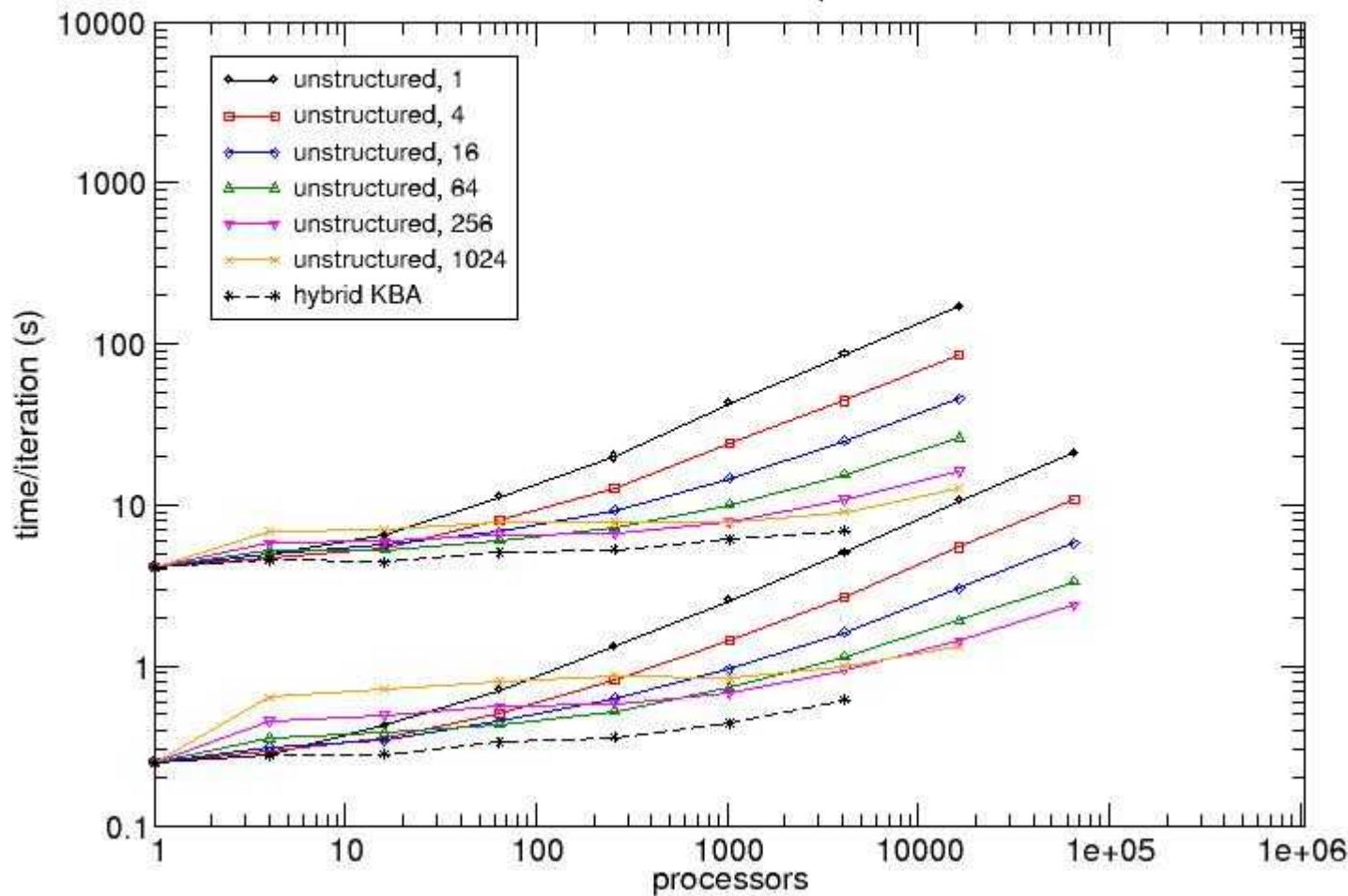
Sceptre first-order weak parallel scaling, Cielo

S2, P0, 5 iterations, NxN quad4 meshes



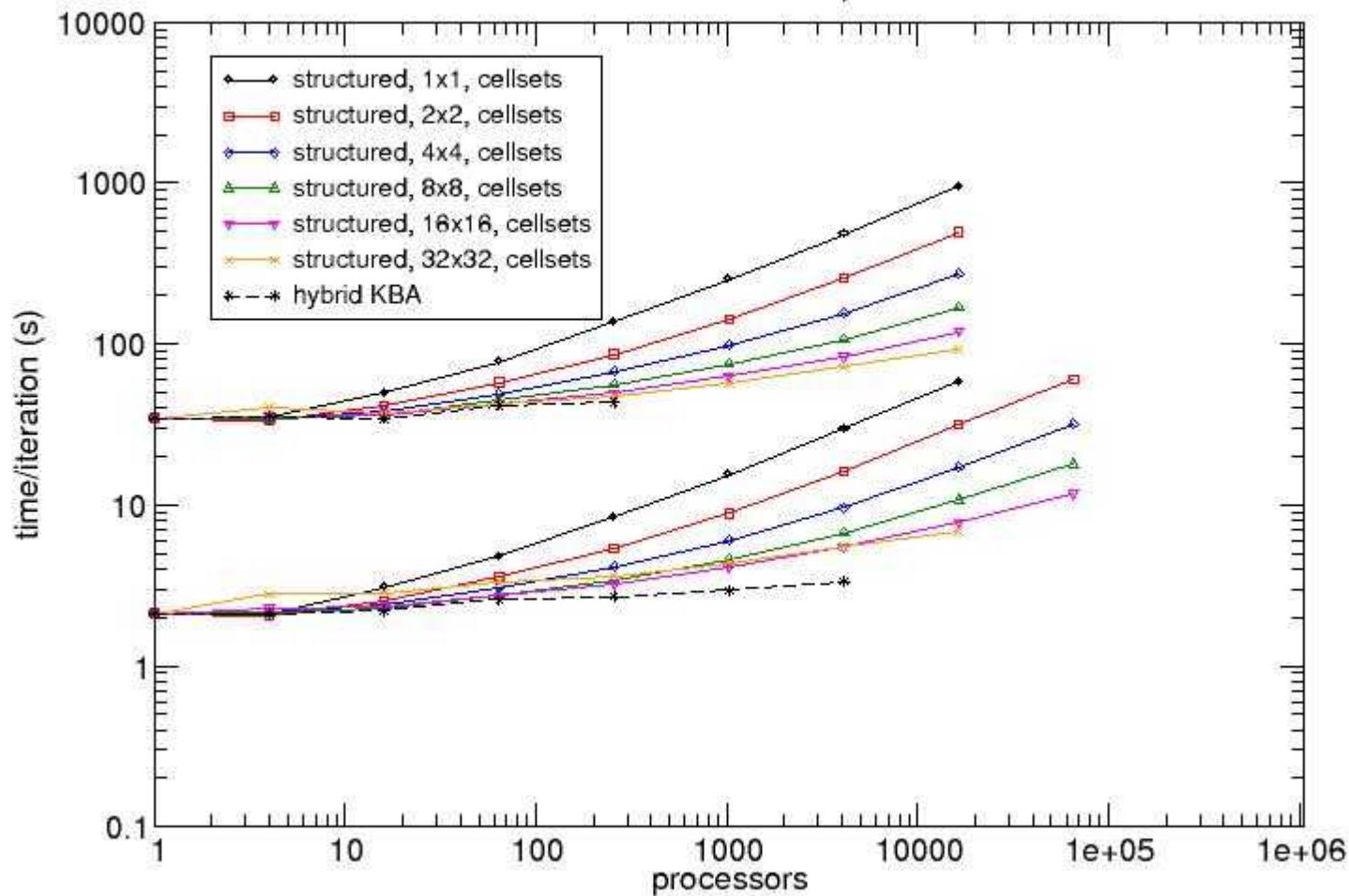
Sceptre first-order weak parallel scaling, Cielo

S2, P0, 5 iterations, NxN quad4 meshes



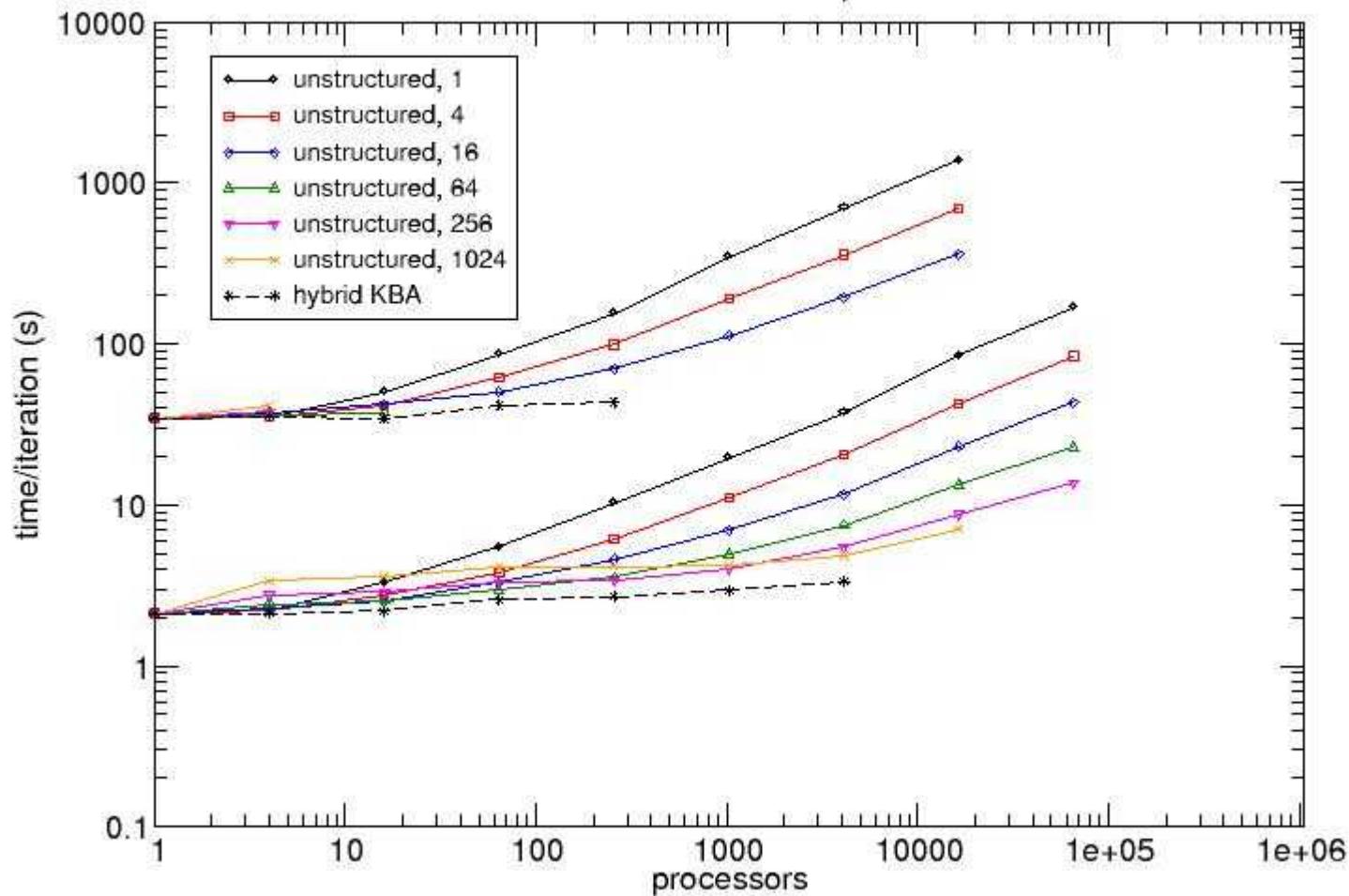
Sceptre first-order weak parallel scaling, Cielo

S8-S4, P0, 5 iterations, NxN quad4 meshes



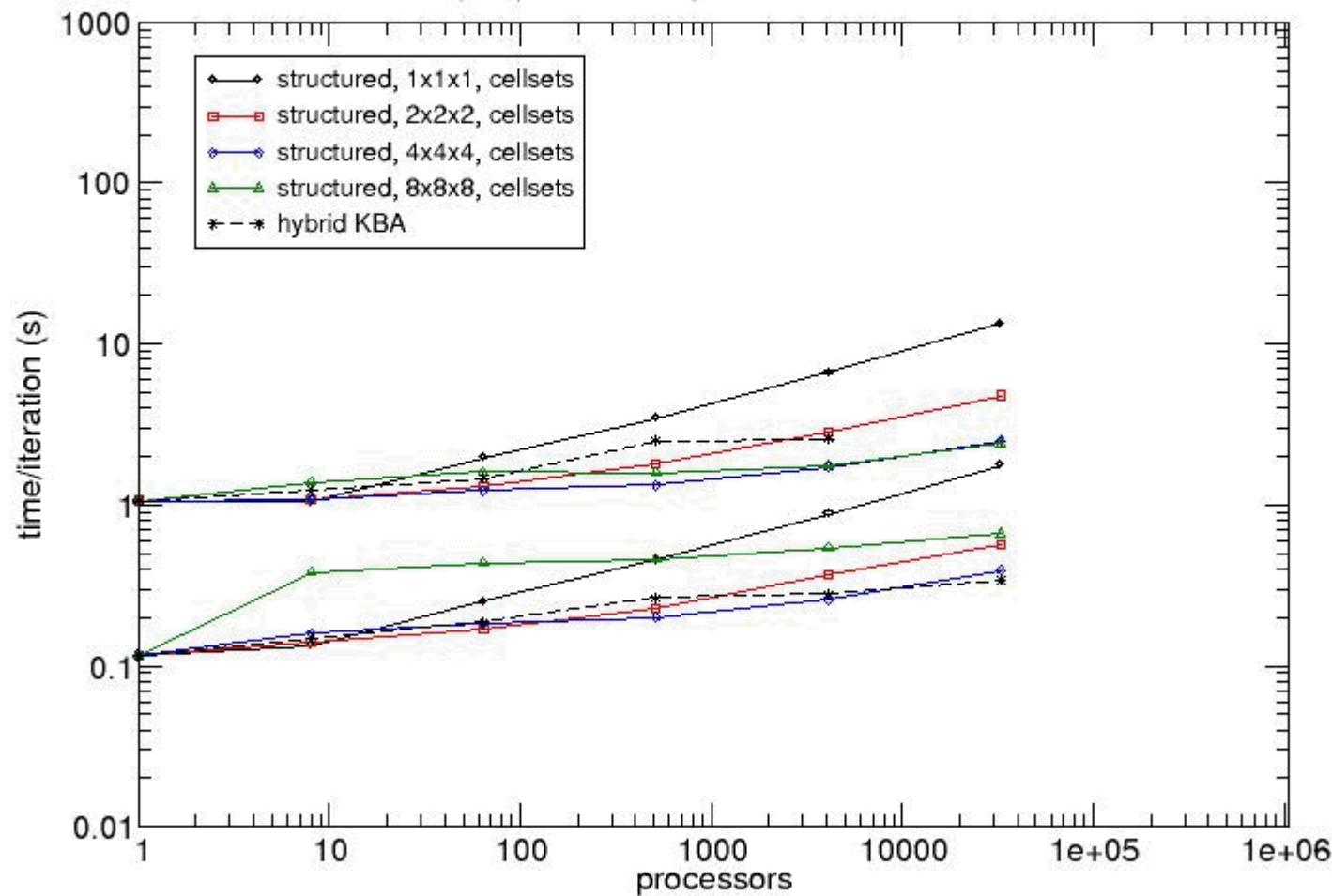
Sceptre first-order weak parallel scaling, Cielo

S8-S4, P0, 5 iterations, NxN quad4 meshes



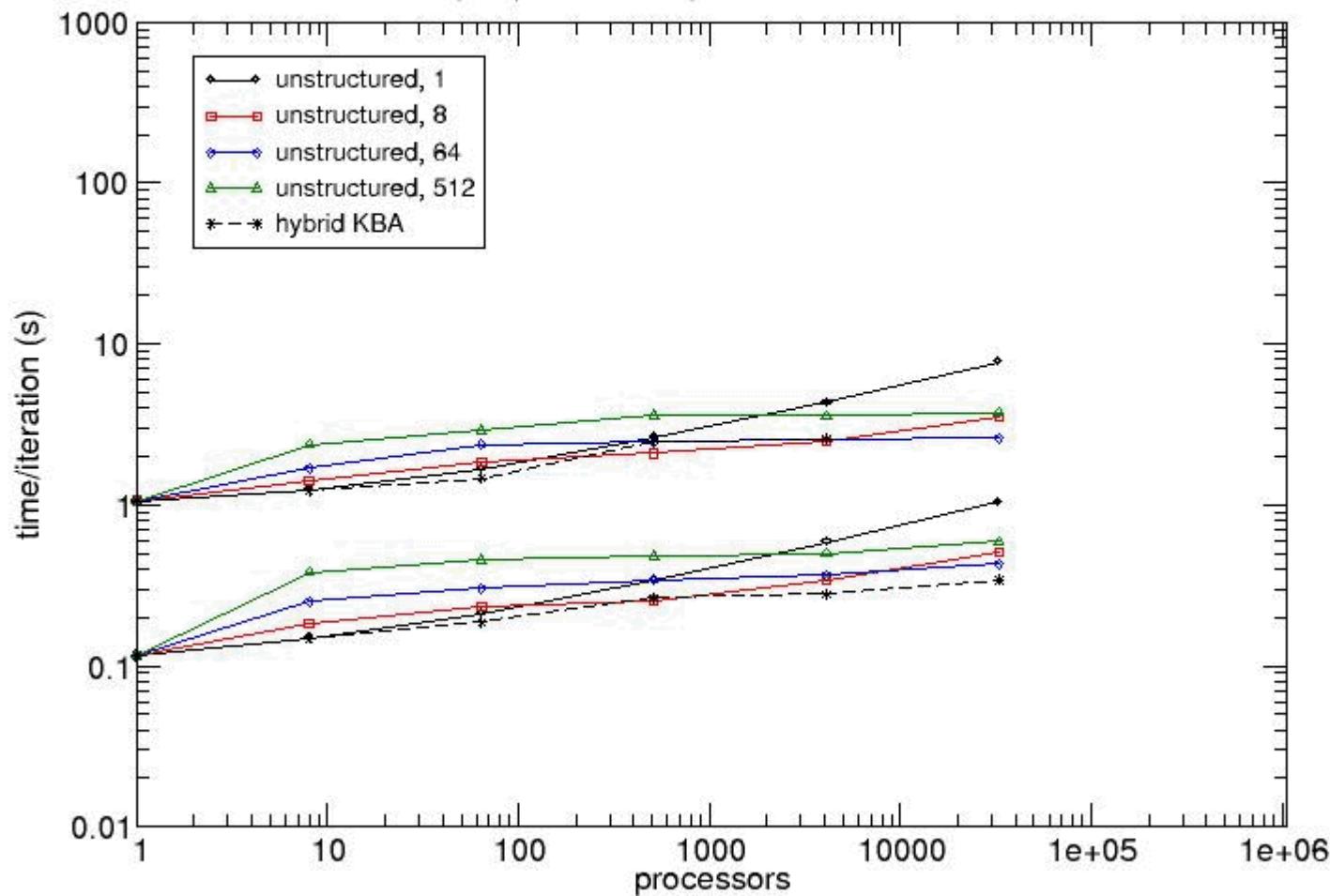
Sceptre first-order weak parallel scaling, Cielo

S2, P0, 5 iterations, NxNxN hex8 meshes



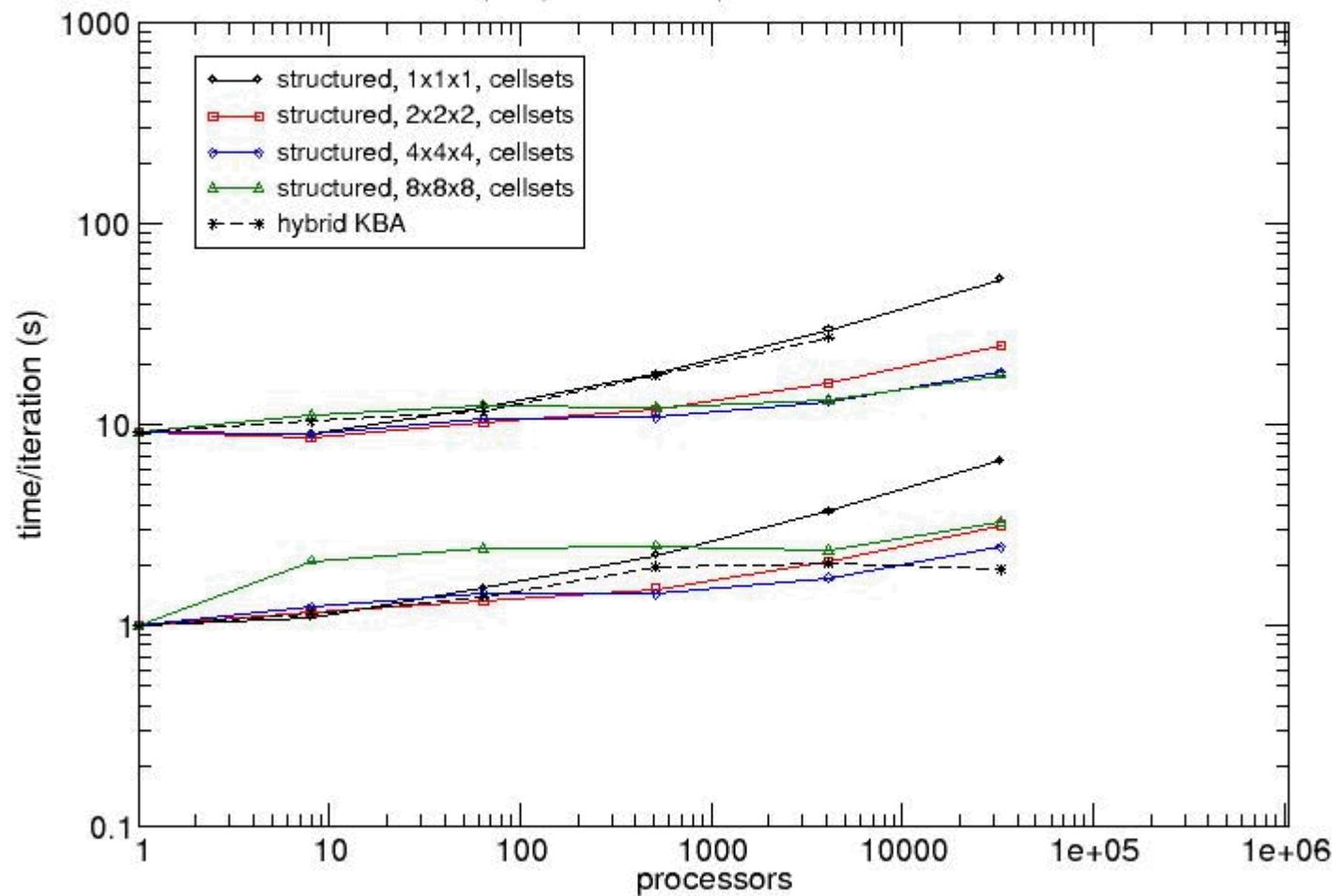
Sceptre first-order weak parallel scaling, Cielo

S2, P0, 5 iterations, NxNxN hex8 meshes



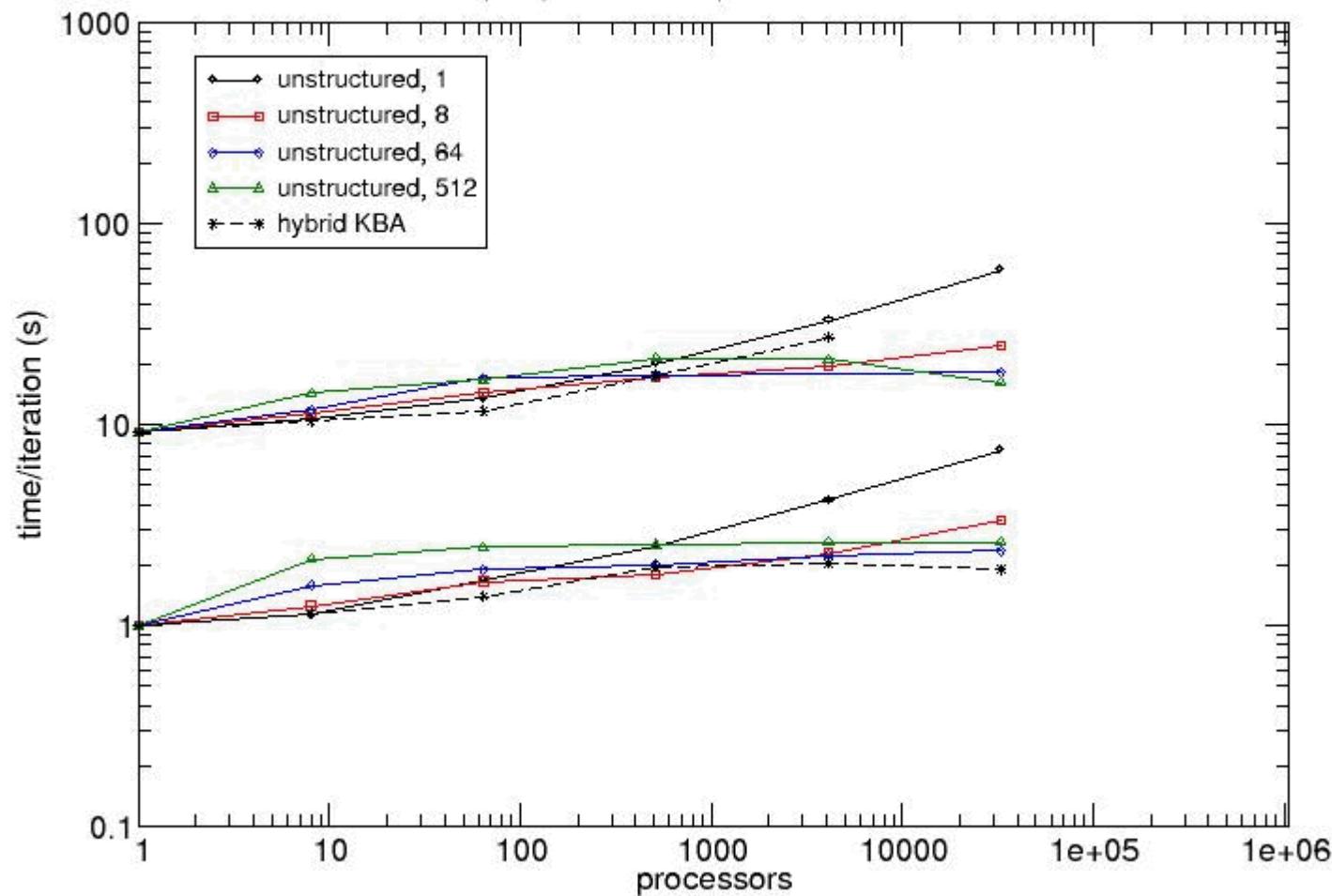
Sceptre first-order weak parallel scaling, Cielo

S8-S4, P0, 5 iterations, NxNxN hex8 meshes



Sceptre first-order weak parallel scaling, Cielo

S8-S4, P0, 5 iterations, NxNxN hex8 meshes



Conclusions

- We have successfully extended the concept of overloading to the general case
- The technique leverages existing tools
- The technique leads to improved scaling for both structured and unstructured algorithms
- More studies are desired to consider other aggregation and scheduling heuristics
- Challenge: performance models for the general case
 - Unstructured vs. structured
 - More general aggregation approaches