A High-Level Approach for Parallelizing Legacy Applications for Multiple Target Platforms

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From a Report from the Council on Competitiveness (Compete.org)

4) Three Systemic Barriers Are Stalling HPC Adoption: Lack of Application Software, <u>Lack</u> of Sufficient Talent, and Cost Constraints

"The problem is price and the need for a dedicated technical person."

"Lack of in-house expertise is a problem for us."

"It could be beneficial to our business if an appropriate return on investment could be realized for the acquisition [of] and training on these tools."







- What do they really want to focus on?
 - Get the science done quickly or spend time in learning low-level details of different parallel programming paradigms?
 - Leverage the investments made in legacy application development or invest in HPC application development from scratch?





Which are the most widely used parallel programming paradigms for High Performance Computing applications? MPI **OpenMP CUDA** Hybrid programming These paradigms can be classified under the category of explicit parallelization.





Traditional Process of Explicit Parallelization







Mastering Multiple Paradigms for Explicit Parallelization is **Often a Challenge for Several Domain-Experts**



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Why is Traditional Explicit Parallelization Using Different Paradigms a Challenge?

- Manual reengineering of legacy applications can be an effortand time-intensive activity even for large and well-funded teams
 - First, need to understand the microarchitectural details of the latest HPC platforms
 - Then, learn about the details of the supported parallel programming paradigms
 - Then, invest time and effort in manually reengineering the code
- What if you do not see good performance at the end of the whole exercise?



- There is a need for a high-level approach that can offer a low-risk way for domain-experts to try HPC
- There is a need for a tool that can help in porting legacy applications to latest HPC platforms



Vision for the Desired High-Level Solution for Explicit Parallelization





Explicit Parallelization at a High-Level



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What is the secret sauce that went into the design and development of our high-level tool ?

- Encapsulated the knowledge of expert parallel programmers inside design templates and rules that are used for source-to-source transformation
- 2. Abstracted the commonly seen standard and nonstandard steps involved in explicit parallelization
- 3. Adopted the user-guided approach instead of 100% automation





Standard and Non-Standard Steps for Parallelization that are Repeatable

- Examples of standard steps in developing an MPI application (common in all MPI programs)
 - Every MPI program has #include "mpi.h"
 - Every MPI program has MPI_Init and MPI_Finalize
 function calls
- Non-standard steps in developing an MPI application
 - for-loop parallelization, data distribution, mapping of tasks to processes, and orchestration of exchange of messages
 - Steps for splitting the work in a for-loop amongst all the processes in MPI_COMM_WORLD are standard for a given load-balancing scheme





Which Interface is the right one?

- Command-Line Interface or the Interactive Parallelization Tool (IPT)
 - Most light-weight
 - Convenient for small applications and low barrier to adoption
- Graphical User Interface or GUI
 - Hides the transformation details from the end-user (those that they really do not want to know about) but is heavier than IPT for remote usage
 - Convenient for small applications and low barrier to adoption
- Domain-Specific Language or Hi-Pal interface
 - Convenient for large applications with repeated patterns and cross-cutting concerns but involves a small learning curve





Parallel Programming Via IPT

- Our Interactive Parallelization Tool (IPT) can be used for transforming legacy C/C++ programs into multiple parallel variants
 - Support for Fortran applications will be added in future
- IPT can be used to teach and learn different parallel programming paradigms through comparison and demonstration
 - without worrying about the low-level details related to the syntax and semantics of different paradigms
- IPT can help in porting legacy applications to latest architectures
- IPT shortens the application development cycle and hence can quickly show the impact of the design choices on performance
 - impact of static load balancing versus dynamic load-balancing
 - impact of choosing MPI only versus choosing hybrid programming





IPT Architecture



CUDA

A Small Parallelization Exercise

```
1. //other code
2. NTIMES = atoi(argv[3]);
3. a = allocMatrix<double>(a, M, N);
4. b = allocMatrix<double>(b, M, N);
5. f = allocMatrix<double>(f, M, N);
6. start = 0;
7. //other code
8. printMatrix<double>(a, M, N);
9. t1 = qettime();
10. for (k = start; k < NTIMES \&\& norm >= tolerance; k++) {
11. b = compute(a, f, b, M, N);
12. ptr = a;
13. a = b;
14. b = ptr;
                                          Code snippet of serial
15. norm = normdiff(b, a, M, N);
                                          Poisson Solver Code
16. }
17. t^2 = qettime(); //other code
```











Generated MPI Code for the Exercise Ó X 2:stampede.tacc.utexas.edu - default - SSH Secure Shell File Edit View Window Help 🔲 🎒 🛕 🔳 🎉 🖻 🛍 💼 👫 🖓 📁 🦓 🤣 🌾 🛐 Quick Connect 🚞 Profiles t1 = qettime(); for (k = 0; (k < NTIMES) && (norm >= tolerance); k++) { b = compute(a,f,b,M,N); b = exchange<double
(b,M+2,N+2,P1,Q1,p1,q1,comm2d, rowcomm, colcomm, diag1c) ----- The template for exchanging omm, diaq2comm); otr = a: the data between the ghost a = b: cells of submatrices on b = ptr; different processes rose norm0 = normdiff(b,a,M,N); MPI Allreduce(&norm,&rose_norm0,1,MPI_DOUBLE,MPI_SUM,MPI_COMM_WORLD); } ---- MPI Reduce function call norm = rose norm0; t2 = gettime(); Setting the value of a variable 97.1 87% to the reduced value SSH2 - aes128-cbc - hmac-md5 - nc 80x14 Connected to stampede.tacc.utexas.edu

Exchange Template

ø 2:stampede.tacc.utexas.edu - default - SSH Secure Sh File Edit View Window Help 🖻 🖻 💾 🎮 🙇 🎾 🧠 餐 🌾 8 4 1 <u>, 196</u> 🛃 Quick Connect 🚞 Profiles template <typename T> T** exchange(T** data, int nrows, int ncols, int P, int Q, int p, int q, MPI Comm comm2d, MPI Comm rowcomm, MPI Comm colcomm) { MPI Aint sizeoftype; MPI Datatype datatype, temptype, vectype; The generated code will have the call to the int next, prev, down, up; MPI Request req[8]; exchange template inserted. The exchange MPI Status status[8]; T tempvar = (T)NULL; template has the code for exchanging the datatype = get_mpi_type(tempvar); data amongst the ghost cells in a stencil-// Create datatype for the recutype MPI Type vector(nrows-2, 1, ncols, datatype, & based code. The generated code is readable MPI Type extent(datatype, &sizeoftype) ; int $blens[2] = \{1, 1\};$ and is easy to understand as comments are MPI_Aint displ[2] = {0, sizeoftype}; MPI_Datatype types[2] = {temptype, MPI_UB}; MPI_Type_struct (2, blens, displ, types, &vect inserted wherever necessary. MPI Type commit(&vectype); MPI Cart shift(rowcomm, 0, -1, &prev, &next); MPI Cart shift(colcomm, 0, -1, &down, &up); // printf("[%d,%d]: next=%d prev=%d down=%d up=%d\n", p, q, next, prev, down, up); // send and receive the boundary rows MPI Irecv(&data[0][1], ncols-2, datatype, up, 0, colcomm, &req[0]); MPI Irecv(&data[nrows-1][1], ncols-2, datatype, down, 0, colcomm, &req[1]); MPI Isend(&data[1][1], ncols-2, datatype, up, 0, colcomm, &req[2]); MPI Isend(&data[nrows-2][1], ncols-2, datatype, down, 0, colcomm, &req[3]); MPI_Irecv(&data[1][0], 1, vectype, next, 0, rowcomm, &req[4]); MPI Irecv(&data[1][ncols-1], 1, vectupe, prev, 0, rowcomm, &req[5]); MPI Isend(&data[1][1], 1, vectype, next, 0, rowcomm, &req[6]);

Benefits of IPT

- In how much time can you manually parallelize the Poisson Solver program using MPI?
 - IPT can help you in parallelizing this code in approximately 5 minutes given that you know the high-level concepts related to parallel programming, and are already familiar with IPT
 - IPT inserted about 357 lines of code in the serial version of the code in order to develop an MPI version
- In how much time can you learn a new parallel programming paradigm and use the knowledge gained in porting legacy application to a new platform manually?
 - IPT significantly reduces the time-to-solution
- The usability study to quantify the benefits related to IPT is pending





Results- Stencil-Based Pattern-16 MPI Processes



MPI - Comparison of Serial and Parallel Code Runtimes





Results- Stencil-Based Pattern-16 OMP Threads



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OpenMP - Comparison of Serial and Parallel Code Runtimes

Results – Stencil-Based Computations - CUDA



CUDA - Comparison of Serial and Parallel Code Runtimes





Run-time Comparison - Circuit Satisfiability Problem







Graphical User Interface (Demo)

O O X Demo:Generator Tool					
File Selection Help					
Programming Model	int solution num:		Target Source		
⊖ CUDA	int value;		a get som es		
O MPI	double wtime1;				
OnenMP	couple wilmez;				
© Openini	timestamp(); printf("\n");				
	ihi = 1;				
	tor $(1 = 1; 1 \le n; 1++)$ { ihi = (ihi * 2);				
)				
	solution_num = 0; wtime1 = gettime0;				
Languages	for $(i = 0; i < ini; i + +)$ {				
• c	i4_to_bvec(i,n,bvec);				
○ C++	if (value = 1) {				
O Fortran	solution_num = (solution_num + 1);				
	for $(j = 0; j < n; j++)$				
	printf("%d",bvec[j]);				
	print("\n"):				
	1				
	<pre>wtime2 = gettime0;</pre>				
Posta o ma	printf("\n");				
	printf(" Number of solutions found was %d\n", solution_num); printf(" Elapsed wall clock time (seconds) %f)n" (witime? - witime1));				
	print('Lipsea nan containe (seconds) a (n),(minez 'minez),,				
CUDACDouble-nested-for-loop	timestamp();				
• OpenMP:Outer-for-loop, reduce	return 0;				
O Stencil	þ				
O Pipeline	int circuit_value(int n,int bvec[])				
	(instantion				
	value = ((((((((((((((((((((((((((((())) a ([1]!=			
	return value;	•			
	Edit Save Cancel Run	1.1.1	Edit Save Cancel		
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Providing Specifications Through Hi-PaL (1)

Parallel section begins <hook type> (<hook pattern>)
mapping is <mapping type> {
 <Hi-PaL API for specifying the operation> <hook> &&
 in function (<function name>)

General Structure of Hi-PaL Code to Generate MPI Code

OMP_Parallel {

<Hi-PaL API for specifying the operation> && schedule is
<schedule type> <hook> && in function (<function name>)

General Structure of Hi-PaL Code to Generate OpenMP Code





Providing Specifications Through Hi-PaL (2)

A set of Hi-PaL API has been developed for precisely capturing the end-users' specifications at a high-level

Hi-PaL API	Description	
<pre>ParExchange2DArrayInt(<array name="">, <num of="" rows="">, <num columns="" of="">)</num></num></array></pre>	Exchange neighboring values in stencil-based computations	
<pre>Parallelize_For_Loop where (<for_init_stmt>; <condition>; <stride>)</stride></condition></for_init_stmt></pre>	Parallelize for-loop with matching condition, stride and initialization statement	
ReduceSumInt(<variable name>)</variable 	MPI_Reduce with MPI_SUM operation or OpenMP reduction clause with '+' operator; reduced variable is of type integer	





Parallelizing Poisson Solver (1)

```
1. //other code
2. NTIMES = atoi(argv[3]);
3. a = allocMatrix<double>(a, M, N);
4. b = allocMatrix<double>(b, M, N);
5. f = allocMatrix<double>(f, M, N);
6. start = 0;
7. //other code
8. printMatrix<double>(a, M, N);
9. t1 = qettime();
10. for (k = start; k < NTIMES \&\& norm >= tolerance; k++) {
11. b = compute(a, f, b, M, N);
12. ptr = a;
13. a = b;
14. b = ptr;
15. norm = normdiff(b, a, M, N);
16. }
17. t_2 = qettime(); //other code
```

Code snippet of serial Poisson Solver Code





Parallelizing Poisson Solver (2)

1. Parallel section begins after ("NTIMES = atoi(argv[3]);")
mapping is Linear{

- 2. ParExchange2DArrayDouble (a, M, N) before statement
 ("printMatrix<double>(a, M, N);")
 - && in function ("main");
- 3. ParExchange2DArrayDouble (b, M, N) before statement
 ("printMatrix<double>(a, M, N);")
 - && in function ("main");
- 4. ParExchange2DArrayDouble (b, M, N) after statement
 ("b=compute(a, f, b, M, N);") && in function ("main");
- 5. AllReduceSumInt(norm) after statement
 ("norm = normdiff(b, a, M, N);") && in function ("main")
 6. }

Hi-PaL Code to Generate MPI Code for Poisson Solver





Generated MPI Code for Poisson Solver (1)

1. //other code

- 2. NTIMES = atoi(argv[3]);
- 3. MPI_Init(NULL, NULL);
- 4. MPI_Comm_size(MPI_COMM_WORLD, &size_Fraspa);
- 5. MPI_Comm_rank(MPI_COMM_WORLD, &rank_Fraspa);
- 6. create_2dgrid(MPI_COMM_WORLD, &comm2d_Fraspa,...);
- 7. create_diagcomm(MPI_COMM_WORLD, size_Fraspa, ...);
- 8. rowmap_Fraspa.init(M, P_Fraspa, p_Fraspa);
- 9. colmap_Fraspa.init(N, Q_Fraspa, q_Fraspa);
- 10. myrows_Fraspa = rowmap_Fraspa.getMyCount();
- 11. mycols_Fraspa = colmap_Fraspa.getMyCount();
- 12. $M_{\text{Fraspa}} = M;$
- 13. $N_Fraspa = N;$
- 14. M = myrows_Fraspa;
- 15. N = mycols_Fraspa;
- 16. a = allocMatrix<double>(a, M, N);





Generated MPI Code for Poisson Solver (2)

```
17. b = allocMatrix<double>(b, M, N);
18. f = allocMatrix<double>(f, M, N);
19. start = 0;
20. //other code
21. a = exchange<double>(a, myrows Fraspa + 2, ...);
22. b = exchange<double>(b, myrows Fraspa + 2, ...);
23. printMatrix<double>(a, M, N);
24. t1 = MPI Wtime();
25. for (k = start; k < NTIMES \&\& norm >= tolerance; k++) {
26. b = compute(a, f, b, M, N);
27. b = exchange<double>(b, myrows Fraspa + 2, ...);
28. ptr = a;
29. a = b;
30. b = ptr;
31. norm = normdiff(b, a, M, N);
32. MPI Allreduce(&norm, &norm Fraspa, 1, MPI INT, MPI SUM, ...);
33. norm = norm Fraspa;
34. }
                               31
36. //other code
```

Results: Poisson Solver (Hi-PaL based MPI)







Results: Genetic Algorithm for Content Based Image Retrieval (Hi-PaL based MPI & OpenMP)







Results: Seismic Tomography Code (**GUI-based CUDA**)







Results: Circuit Satisfiability Code (GUI-based OpenMP + Offload)







Summary of Features & Benefits

- Enhances the productivity of the end-users in terms of the reduction in the time and effort
 - reduction in manual effort by over 90% while ensuring that the performance of the generated parallel code is within 5% of the sample hand-written parallel code
- Leverages the knowledge of expert parallel programmers
- Separates the sequential and parallel programming concerns while preserving the existing version of sequential applications





Ongoing and Future Work

- Code cleanup and preparation for the public release
- Support for handling irregular meshes will be added in future
- Support for Fortran code generation will be added in future as well
- Support for generating hybrid applications is available with Hi-PaL interface. The command-line interface and GUI will be extended to support hybrid code generation as well





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Questions or Comments?





Separation of Sequential and Parallel Concerns





