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 Х 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)

Image: Selection Help Image: Selection Help					
O CUDA	int value;		Car gar sen ca		
O MPI	double wtime 1;				
OpenMP	double wtime2; printf("\n");				
© openni	timestamp();				
	printf("\n");				
	ihi = 1; for (i = 1; i <= n; i++) {				
	ihi = (ihi * 2);				
	}				
	solution_num = 0; wtime1 = gettime();				
Languages	for $(i = 0; i < ihi, i++)$ {				
• c	i4_to_bvec(i,n,bvec); value = circuit_value(n,bvec);				
0 C++	if (value = 1) {				
O Fortran	solution_num = (solution_num + 1);	=			
Fordan	printf(" %2d %10d: ",solution_num,i); for (j = 0; j < n; j++) {				
	printf(":%d", bvec[j]);				
	3				
	printf("\n");				
	3				
	wtime2 = gettime();				
Patterns	printf("\n"); printf(" Number of solutions found was %d\n", solution_num);				
○ CUDA:Outer-for-loop	printf(" Elapsed wall clock time (seconds) %f\n", (wtime2 - wtime1));				
CUDA:Double-nested-for-loop	printf('\n'');				
OpenMP:Outer-for-loop, reduce	timestamp(); printf("\n");				
	return 0;				
O Stencil)				
O Pipeline	int circuit_value(int n, int bvec[])				
	K				
	int value; value = ((((((((((((((((((((((((((((((((())) term ())))))))))))))))))))))))))))))))))))	111= 0			
	return value;	-		•	
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	Edit Save Cancel Run		Edit Save Cancel	1	
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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





