XcalableACC: Extension of XcalableMP using OpenACC for Accelerator Clusters

Motivation

In order to develop an application on accelerator clusters, a programmer selects a combination of MPI and one of the available accelerator programming models, such as OpenACC, OpenCL or CUDA. In particular, OpenACC provides good productivity and high performance. However, MPI programming is often difficult because programmers need to distribute, transfer, and reduce data by using primitive MPI functions.

XcalableACC

Based on studies of XMP and OpenACC, we believe that a directive-based language is effective for also accelerator clusters. Thus, we propose a new programming model called *XcalableACC (XACC)*, which is an XMP extension that uses OpenACC. In XACC, programmers can use both XMP and OpenACC directives seamlessly. Therefore, XACC allows programmers to easily develop applications on accelerator clusters. Additionally, XACC also provides directives to transfer data among accelerators efficiently. **"Example 1"** on the right side shows an example of XACC.

Extension of OpenACC for multiple accelerators

In the OpenACC specification 2.0a, it is difficult for a single process to use multiple accelerators. For example, programmers must specify an accelerator number using the **acc_set_device_num()** before each statement, calculate indexes for loop statement on each GPU, and move data among accelerators manually. To reduce the programming cost, we consider an extension of OpenACC to use multiple accelerators. We believe that this extension is also useful for XACC. **"Example 2"** on the right side shows an example of the extension of OpenACC.

Definition and synchronization of a halo region

To develop stencil applications easily, XACC provides shadow and reflect directives. These directives can be used to define a halo region for a distributed array and synchronize it among nodes. In order to transfer data among accelerator memories efficiently, the

Example 1: Loop parallelization



Example 2: Usage of multiple accelerators



Evaluation of the HIMENO benchmark

The HIMENO Benchmark is a stencil application benchmark that evaluates the performance of incompressible fluid analysis code.

<pre>float p[MIMAX][MJMAX][MKMAX]; #pragma xmp shadow p[1:1][1:1][0]</pre>	Define a halo region
<pre>#pragma xmp reflect (p) acc</pre>	Synchronize a halo region
<pre>#pragma xmp loop (k,j,i) on t(k,j,i) #pragma acc parallel loop collapse(3)</pre>	Parallelize a loop statement
<pre>for(i=1 ; i<mimax ++i)="" ++j){<="" ;="" for(j="1" j<mjmax="" pre=""></mimax></pre>	
for(k=1 ; k <mkmax ++k){<br="" ;="">S0 = p[i+1][j][k] *;</mkmax>	

Productivity

	Total	Directive		Except for
		XcalableMP	OpenACC	directives
XcalableACC	204	22	9	173
MPI + OpenACC	325	-	15	310

The source lines of the VACC HIMENIC Dependements is less than that of

Omni compiler supports *Tightly Coupled Accelerators(TCA)* in addition to the combination of CUDA and MPI, and GPUDirect RDMA over InfiniBand. TCA is a communication architecture based on PCI Express (PCIe) technology. Computation hosts with TCA network are connected with each other using PCIe external cables, which permits direct, low latency data transfers among accelerator memories because they do not need host memory copies, a protocol conversion, and MPI software stack. For more details of TCA,

please go to "Center for Computational Sciences, University of Tsukuba" (#3215 booth).



the MPI+OpenACC HIMENO Benchmark is less than that of XACC HIMENO benchmark, a programmer only adds XMP and OpenACC directives into the sequential Himeno Benchmark.

Performance



Computer simulations create the future