OP2 C++ User's Manual

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1 Introduction

OP2 is a high-level framework with associated libraries and preprocessors to generate parallel executables for applications on unstructured grids. This document describes the C++ API, but FORTRAN 90 is also supported with a very similar API.

The key concept behind OP2 is that unstructured grids can be described by a number of sets. Depending on the application, these sets might be of nodes, edges, faces, cells of a variety of types, far-field boundary nodes, wall boundary faces, etc. Associated with these are data (e.g. coordinate data at nodes) and mappings to other sets (e.g. edge mapping to the two nodes at each end of the edge). All of the numerically-intensive operations can then be described as a loop over all members of a set, carrying out some operations on data associated directly with the set or with another set through a mapping.

OP2 makes the important restriction that the order in which the function is applied to the members of the set must not affect the final result. This allows the parallel implementation to choose its own ordering to achieve maximum parallel efficiency. Two other restrictions are that the sets and maps are static (i.e. they do not change) and the operands in the set operations are not referenced through a double level of mapping indirection (i.e. through a mapping to another set which in turn uses another mapping to data in a third set).

OP2 currently enables users to write a single program which can be built into three different executables for different single-node platforms:

- single-threaded on a CPU
- parallelised using CUDA for NVIDIA GPUs
- multi-threaded using OpenMP for multicore x86 systems

In the future, there will also be support for AVX vectorisation for x86 CPUs, and OpenCL for both CPUs and GPUS.

In addition to this, there is support for distributed-memory MPI parallelisation in combination with any of the above. The user can either use OP2's parallel file I/O capabilities for HDF5 files with a specified structure, or perform their own parallel file I/O using custom MPI code.

2 Overview

A computational project can be viewed as involving three steps:

- writing the program
- debugging the program, often using a small testcase
- running the program on increasingly large applications

With OP2 we want to simplify the first two tasks, while providing as much performance as possible for the third.

To achieve the high performance for large applications, a preprocessor is needed to generate the CUDA code for GPUs or OpenMP code for multicore x86 systems. However, to keep the initial development simple, the single-threaded executable does not use any special tools; the user's main code is simply linked to a set of library routines, most of which do little more than error-checking to assist the debugging process by checking the correctness of the user's program. Note that this single-threaded version will not execute efficiently. The preprocessor is needed to generate efficient OpenMP code for x86 systems.

Figure 1 shows the build process for a single thread CPU executable. The user's main program (in this case jac.cpp) uses the OP header file op_seq.h and is linked to the appropriate OP libraries using g++, perhaps controlled by a Makefile.

Figure 2 shows the build process for the corresponding CUDA executable. The preprocessor parses the user's main program and produces a modified main program and a CUDA file which includes a separate file for each of the kernel functions. These are then compiled and linked to the OP libraries using g++ and the NVIDIA CUDA compiler nvcc, again perhaps controlled by a Makefile.

Figure 3 shows the OpenMP build process which is very similar to the CUDA process except that it uses *.cpp files produced by the preprocessor instead of *.cu files.

In looking at the API specification, users may think it is a little verbose in places. e.g. users have to re-supply information about the datatype of the datasets being used in a parallel loop. This is a deliberate choice to simplify the task of the preprocessor, and therefore hopefully reduce the chance for errors. It is also motivated by the thought that "programming is easy; it's debugging which is difficult". i.e. writing code isn't time-consuming, it's correcting it which takes the time. Therefore, it's not unreasonable to ask the programmer to supply redundant information, but be assured that the preprocessor or library will check that all redundant information is self-consistent. If you declare a dataset as being of type OP_DOUBLE and later say that it is of type OP_FLOAT this will be flagged up as an error at run-time.

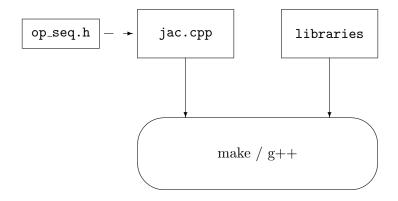


Figure 1: Sequential code build process

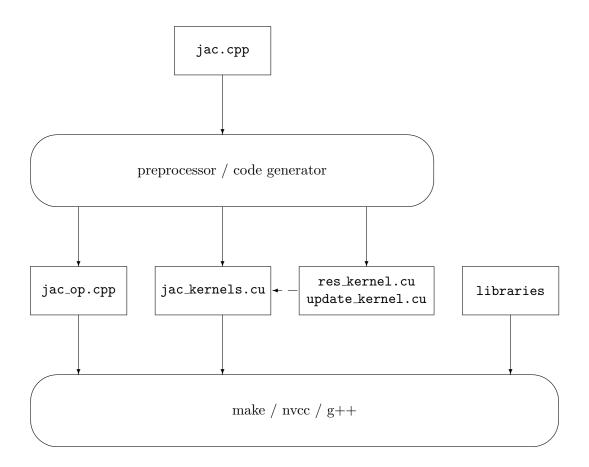


Figure 2: CUDA code build process

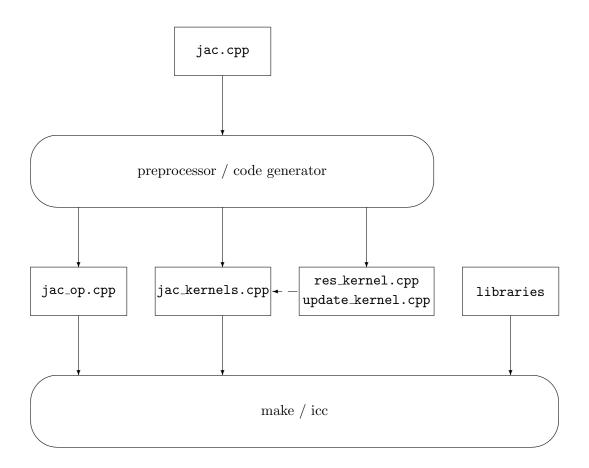


Figure 3: OpenMP code build process

$3 \quad OP2 C++ API$

3.1 Initialisation and termination routines

void op_init(int argc, char **argv, int diags_level)

This routine must be called before all other OP routines.

argc, argv the usual command line arguments

diags_level an integer which defines the level of debugging diagnostics and reporting

to be performed;

0 - none;

1 - error-checking;

2 – info on plan construction;

3 – report execution of parallel loops;

4 – report use of old plans;

7 – report positive checks in op_plan_check;

void op_exit()

This routine must be called last to cleanly terminate the OP computation.

op_set op_decl_set(int size, char *name)

This routine defines a set, and returns a set ID.

size number of elements in the set

name a name used for output diagnostics

op_map op_decl_map(op_set from, op_set to, int dim, int *imap, char *name)

This routine defines a mapping from one set to another, and returns a map ID.

from set pointed from

to set pointed to

dim number of mappings per element

imap input mapping table

name a name used for output diagnostics

void op_decl_const(int dim, char *type, T *dat, char *name)

This routine declares constant data with global scope to be used in user's kernel functions. Note: in sequential version, it is the user's responsibility to define the appropriate variable with global scope.

dim dimension of data (i.e. array size)

for maximum efficiency, this should be a literal constant (i.e. a number

not a variable)

type datatype, either intrinsic ("float", "double", "int", "uint", "ll", "ull" or

"bool") or user-defined

dat input data of type T (checked for consistency with type at run-time)

name global name to be used in user's kernel functions;

a scalar variable if dim=1, otherwise an array of size dim

op_dat op_decl_dat(op_set set, int dim, char *type, T *data, char *name)

This routine defines a dataset, and returns a dataset ID.

 $\mathtt{set} \qquad \qquad \mathrm{set}$

dim dimension of dataset (number of items per set element)

at present this must be a literal constant (i.e. a number not a variable); this restriction will be removed in the future but a literal constant will

remain more efficient

type datatype, either intrinsic or user-defined

data input data of type T (checked for consistency with type at run-time)

name a name used for output diagnostics

void op_fetch_data(op_dat dat)

This routine transfers data from the GPU back to the CPU.

dat OP dataset ID – data is put back into original input array

void op_diagnostic_output()

This routine prints out various useful bits of diagnostic info about sets, mappings and datasets

3.2 Parallel loop syntax

A parallel loop with N arguments has the following syntax:

void op_par_loop(void (*kernel)(...), char *name, op_set set, op_arg arg1, op_arg arg2, ..., op_arg argN)

kernel user's kernel function with N arguments

(this is only used for the single-threaded CPU build)

name of kernel function, used for output diagnostics

set OP set ID

args arguments

The **op_arg** arguments in **op_par_loop** are provided by one of the following routines, one for global constants and reductions, and the other for OP2 datasets. In the future there will be a third one for sparse matrices to support the needs of finite element calculations.

op_arg op_arg_gbl(T *data, int dim, char *typ, op_access acc)

data data array

dim array dimension

typ datatype (redundant info, checked at run-time for consistency)

acc access type:

OP_READ: read-only

OP_INC: global reduction to compute a sum

OP_MAX: global reduction to compute a maximum OP_MIN: global reduction to compute a minimum

op_arg op_arg_dat(op_dat dat, int idx, op_map map, int dim, char *typ, op_access acc)

dat OP dataset ID

idx index of mapping to be used $(-1 \equiv \text{no mapping indirection})$

map OP mapping ID (OP_ID for identity mapping, i.e. no mapping indirection)

dim dataset dimension (redundant info, checked at run-time for consistency)

at present this must be a literal constant (i.e. a number not a variable); this restriction will be removed in the future but a literal constant will

remain more efficient

typ dataset datatype (redundant info, checked at run-time for consistency)

acc access type:

OP_READ: read-only

OP_WRITE: write-only, but without potential data conflict OP_RW: read and write, but without potential data conflict OP_INC: increment, or global reduction to compute a sum

The restriction that OP_WRITE and OP_RW access must not have any potential data conflict means that two different elements of the set cannot through a mapping indirection reference the same elements of the dataset.

Furthermore, with OP_WRITE the user's kernel function must set the value of all DIM components of the dataset. If the user's kernel function does not set all of them, the access should be specified to be OP_RW since the kernel function needs to read in the old values of the components which are not being modified.

op_arg op_arg_dat_opt(op_dat dat, int idx, op_map map, int dim, char *typ, op_access acc, int flag)

This is the same as op_arg_op_arg_dat except for an extra variable flag; the argument is only actually used if flag has a non-zero value. This routine is required for large application codes (such as HYDRA) which have lots of different features turned on and off by logical flags.

Note that if the user's kernel needs to know the value of flag then this must be passed as an additional op_arg_gbl argument.

3.3 MPI message-passing using HDF5 files

HDF5 has become the *de facto* standard format for parallel file I/O, with various other standards like CGNS layered on top. To make it as easy as possible for users to develop distributed-memory OP2 applications, we provide alternatives to some of the OP2 routines in which the data is read by OP2 from an HDF5 file, instead of being supplied by the user:

- op_decl_set_hdf5: similar to op_decl_set but with size replaced by char *file which defines the HDF5 file from which size is read using keyword name
- op_decl_map_hdf5: similar to op_decl_map but with imap replaced by char *file from which the mapping table is read using keyword name
- op_decl_dat_hdf5: similar to op_decl_dat but with dat replaced by char *file from which the data is read using keyword name

In addition, there are the following two routines.

T op_get_const_hdf5(int dim, char *type, char *file, char *name)

This routine reads a constant (or constant array) from an HDF5 file; if required, the user must then call **op_decl_const** to declare it to OP2.

dim	dimension of data (i.e. array size)
	for maximum efficiency, this should be a literal constant (i.e. a number not a variable)
type	datatype, either intrinsic ("float", "double", "int", "uint", "ll", "ull" or "bool") or user-defined; checked at run-time for consistency with T
file	name of the HDF5 file
name	global name to be used in user's kernel functions;
	a scalar variable if dim=1, otherwise an array of size dim

void op_partition(char *method)

This routine controls how the various sets are partitioned.

Method A string which declares the partitioning method to be used. (This needs to be updated with details of how to specify and control parmetis and PT-Scotch partitioning.)

void op_print(char *string)

This routine simply prints a message; this is in place of the standard printf function which would print the same message from each MPI process.

Using the above routines, OP2 will take care of everything, reading in all of the sets, mapping and data, partitioning the sets appropriately, renumbering sets as needed, constructing import/export halo lists, etc., and then performing the parallel computation with halo exchange when needed.

Both MPI and single process executables can be generated, depending on the libraries which are linked in.

3.4 MPI message-passing without HDF5 files

Some users will prefer not to use HDF5 files, or at least not to use them in the way prescribed by OP2. To support these users, an application code may do its own file I/O, and then provide the required data to OP2 using the standard routines.

In an MPI application, multiple copies of the same program are executed as separate processes, often on different nodes of a compute cluster. Hence, the OP2 declarations will be invoked on each process. In this case, the behaviour of the OP2 declaration routines is as follows:

- op_decl_set: size is the number of elements of the set which will be provided by this MPI process
- op_decl_map: imap provides the part of the mapping table which corresponds to its share of the from set
- op_decl_dat: dat provides the data which corresponds to its share of set

For example, if an application has 4 processes, 4×10^6 nodes and 16×10^6 edges, then each process might be responsible for providing 10^6 nodes and 4×10^6 edges. Process 0 (the one with MPI rank 0) would be responsible for providing the first 10^6 nodes, process 1 the next 10^6 nodes, and so on, and the same for the edges.

The edge \rightarrow node mapping tables would still contain the same information as in a single process implementation, but process 0 would provide the first 4×10^6 entries, process 1 the next 4×10^6 entries, and so on.

This is effectively using a simple contiguous block partitioning of the datasets, but it is very important to note that this will not be used for the parallel computation. OP2 will re-partition the datasets, re-number the mapping tables as needed (as well as constructing import/export lists for halo data exchange) and will move all data/mappings/datasets to the correct MPI process.

4 Preprocessor

There are two preprocessors for OP2, one developed at Imperial College using ROSE, and the other developed at Oxford using MATLAB.

```
The MATLAB preprocessor is run by the command op2('main')
```

where main.cpp is the user's main program. It produces as output

- a modified main program main_op.cpp which is used for both the CUDA and OpenMP executables;
- for the CUDA executable, a new CUDA file main_kernels.cu which includes one or more files of the form xxx_kernel.cu containing the CUDA implementations of the user's kernel functions;
- for the OpenMP executable, a new C++ file main_kernels.cpp which includes one or more files of the form xxx_kernel.cpp containing the CUDA implementations of the user's kernel functions.

```
If the user's application is split over several files it is run by a command such as op2('main','sub1','sub2','sub3')
```

where sub1.cpp, sub2.cpp, sub3.cpp are the additional input files which will lead to the generation of output files sub1_op.cpp, sub2_op.cpp, sub3_op.cpp in addition to main_op.cpp, main_kernels.cu, main_kernels.cpp and the individual kernel files.

5 Error-checking

At compile-time, there is a check to ensure that CUDA 3.2 or later is used when compiling the CUDA executable; this is because of compiler bugs in previous versions of CUDA.

At run-time, OP2 checks the user-supplied data in various ways:

- checks that a set has a strictly positive number of elements
- checks that a map has legitimate mapping indices, i.e. they map to elements within the range of the target set
- checks that variables have the correct declared type

It would be great to get feedback from users on suggestions for additional error-checking.

6 32-bit and 64-bit CUDA

Section 3.1.6 of the CUDA 3.2 Programming Guide says:

The 64-bit version of **nvcc** compiles device code in 64-bit mode (i.e. pointers are 64-bit). Device code compiled in 64-bit mode is only supported with host code compiled in 64-bit mode.

Similarly, the 32-bit version of nvcc compiles device code in 32-bit mode and device code compiled in 32-bit mode is only supported with host code compiled in 32-bit mode.

The 32-bit version of nvcc can compile device code in 64-bit mode also using the -m64 compiler option.

The 64-bit version of nvcc can compile device code in 32-bit mode also using the -m32 compiler option.

On Windows and Linux systems, there are separate CUDA download files for 32-bit and 64-bit operating systems, so the version of CUDA which is installed matches the operating system. i.e. the 64-bit version is installed on a 64-bit operating system.

Mac OS X can handle both 32-bit and 64-bit executables, and it appears that it is the 32-bit version of nvcc which is installed. Therefore the Makefiles in the OP2 distribution may need the -m64 flag added to NVCCFLAGS to produce 64-bit object code.

The Makefiles in the OP2 distribution assume 64-bit compilation and therefore they link to the 64-bit CUDA runtime libraries in /lib64 within the CUDA toolkit distribution. This will need to be changed to /lib for 32-bit code.

7 User-defined datatypes

If the user defines a new datatype mytype it must be included in a header file along with

• a type-checking routine:

```
inline int type_error(const mytype *,const char *type)
{return strcmp(type,"mytype");}
```

which is used at run-time to check the consistency of the user's type declarations in input arguments.

• a "zero element" declaration of the form:

```
#define ZERO_mytype 0;
```

as well as an appropriate overloaded addition operator if there is any OP_INC access to the datatype. The zero element and overloaded addition have to be such that 0+x=x where x represents any element of the user's datatype and 0 represents the declared zero element.

• an overloaded implementation of the inequality operators < and > if there are any OP_MIN, OP_MAX accesses to the datatype.

In addition, the user must specify the name of the new header file using the environment variable OP_USER_DATATYPES so that this header file is included into the OP2 header file op_datatypes.h.