
ParPyDTK2 Documentation

Release 1.0.0

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Sep 07, 2018

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INTRODUCTION

1.1 Background & Motivation

Multi-physics coupling problems have been one of the popular topics. People try to put different models together and simulate the behaviors in a coupled fashion. With the popularity of *partitioned approach*, i.e. solve different domains with different solvers and couple the interface conditions as those solvers boundary conditions, a robust and accurate interface solution remapping operator is needed. The solution transfer problem on its own is not an easy task, since it involves the following research aspects:

1. **numerical method**, i.e. consistent, conservative, high-order convergence.
2. **geometry and data structure, i.e. efficient and robust treatments of *mesh association*** of two (potentially more) general surfaces that come from different discretization methods (*FEM, FVM, FDM*, etc.) thus having different resolutions.
3. **parallel rendezvous & HPC, i.e. handling migrating meshes that have** different parallel partitions.

Data Transfer Kit-2.0 (DTK2) is a package that is developed at the Oak Ridge National Laboratory. **DTK2** provides parallel solution transfer services with *meshless* (a.k.a. *mesh-free*) methods, which are relatively easy to implement and computational efficient. Particularly, we are interested in its *modified moving least square*¹ method that is an improvement of traditional MLS fitting in terms of robustness on featured geometries.

Mesh-Oriented datABase is an array-based general purpose mesh library with MPI support. Array-based mesh data structure is more efficient in both computational cost and memory usage compared to traditional pointer-based data structures. **MOAB** has been adapted in **DTK2**, so we choose to use it as our mesh database for this work.

In multi-physics coupling, a flexible software framework is must. The fact is: the physics solvers may be implemented in different programming languages or shipped as executable binaries (typically commercial codes), this makes using static languages difficult. Python, on the other hand, can easily glue different languages together and drive executable binaries smoothly. Its built-in reference counting, garbage collection, and pass-by-reference make it as one of the best choices for developing multi-physics coupling frameworks. In addition, MPI is well supported through the **mpi4py** package. **This is the core motivation of this project!**

1.2 About Me

I am a Ph.D. candidate who work with **Dr. Jim Jiao** on high-order numerical methods. This work is for testing our software framework of multi-physics coupling in general, *conjugate heat transfer* (CHT) in particular.

¹ Slattery, S. Hamilton, T. Evans, “A Modified Moving Least Square Algorithm for Solution Transfer on a Spacer Grid Surface”, ANS MC2015 - Joint International Conference on Mathematics and Computation (M&C), Supercomputing in Nuclear Applications (SNA) and the Monte Carlo (MC) Method, Nashville, Tennessee · April 19–23, 2015, on CD-ROM, American Nuclear Society, LaGrange Park, IL (2015).

Note: Please be aware that I may not have time to maintain this package.

INSTALLATION

Installing this package is not a trivial task due to the heavy dependencies. ParPyDTK2 has the following installation requirements:

1. C++11 compiler
2. MPI
3. [MOAB](#)
4. [DTK2](#) and [Trilinos](#)
5. Python ≥ 3.5
6. [mpi4py](#)
7. [NumPy](#)
8. [setuptools](#)

And both [MOAB](#) and [DTK2](#) have their own dependencies.

In addition, to build the documentation, the following packages are needed:

1. Sphinx
2. Doxygen
3. breathe
4. numpdoc

The good news is you can install these easily through `pip`.

2.1 Install MOAB

The [MOAB official README](#) has a very clear description of the installation process. Here we take an excerpt from our [MOAB Docker image build](#):

```
$ git clone --depth=1 https://bitbucket.org/fathomteam/moab.git
$ cd moab
$ autoreconf -fi
$ ./configure \
  --prefix=/usr/local \
  --with-mpi \
  CC=mpicc \
  CXX=mpicxx \
  FC=mpif90 \
```

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```

F77=mpif77 \
--enable-optimize \
--enable-shared=yes \
--with-blas=-lopenblas \
--with-lapack=-lopenblas \
--with-scotch=/usr/lib \
--with-metis=/usr/lib/x86_64-linux-gnu \
--with-eigen3=/usr/include/eigen3 \
--with-x \
--with-cgns \
--with-netcdf \
--with-hdf5=/usr/lib/hdf5-openmpi \
--with-hdf5-ldflags="-L/usr/lib/hdf5-openmpi/lib" \
--enable-ahf=yes \
--enable-tools=yes
$ make && sudo make install

```

Notice that this is for system installation. Install to your preferred locations if you don't have root access. Also, turn off those optional packages if you don't have them, the only MPI and HDF5 are necessary.

Warning: You must build it into a shared object!

Note: If you use Ubuntu >= 17.10, all these packages are officially supported.

2.2 Install DTK2

DTK2 is shipped as a sub-module of [Trilinos](#), so installing [Trilinos](#) is needed. For people who are not familiar with [Trilinos](#), this can be tricky. The following is an excerpt from our [DTK2 Docker image](#) build:

```

$ export TRILINOS_VERSION=12-12-1
$ git clone --depth 1 --branch trilinos-release-${TRILINOS_VERSION}
$ cd Trilinos
$ git clone --depth 1 --branch dtk-2.0 \
  https://github.com/unifem/DataTransferKit.git
$ mkdir build && cd build
$ cmake \
  -DCMAKE_INSTALL_PREFIX:PATH=/usr/local \
  -DCMAKE_BUILD_TYPE:STRING=RELEASE \
  -DCMAKE_VERBOSE_MAKEFILE:BOOL=OFF \
  -DCMAKE_SHARED_LIBS:BOOL=ON \
  -DTPL_ENABLE_MPI:BOOL=ON \
  -DTPL_ENABLE_Boost:BOOL=ON \
  -DBoost_INCLUDE_DIRS:PATH=/usr/include/boost \
  -DTPL_ENABLE_Libmesh:BOOL=OFF \
  -DTPL_ENABLE_MOAB:BOOL=ON \
  -DMOAB_INCLUDE_DIRS=${MOAB_ROOT}/include \
  -DMOAB_LIBRARY_DIRS=${MOAB_ROOT}/lib \
  -DTPL_ENABLE_Netcdf:BOOL=ON \
  -DTPL_ENABLE_BinUtils:BOOL=OFF \
  -DTrilinos_ENABLE_ALL_OPTIONAL_PACKAGES:BOOL=OFF \

```

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```

-DTrilinos_ENABLE_ALL_PACKAGES=OFF \
-DTrilinos_EXTRA_REPOSITORIES="DataTransferKit" \
-DTrilinos_ENABLE_EXPLICIT_INSTANTIATION:BOOL=ON \
-DTrilinos_ASSERT_MISSING_PACKAGES:BOOL=OFF \
-DTrilinos_ENABLE_TESTS:BOOL=OFF \
-DTrilinos_ENABLE_EXAMPLES:BOOL=OFF \
-DTrilinos_ENABLE_CXX11:BOOL=ON \
-DTrilinos_ENABLE_Tpetra:BOOL=ON \
-DTpetra_INST_INT_UNSIGNED_LONG:BOOL=ON \
-DTPL_ENABLE_BLAS:BOOL=ON \
-DTPL_BLAS_LIBRARIES=/usr/lib/x86_64-linux-gnu/libopenblas.so \
-DTPL_ENABLE_LAPACK:BOOL=ON \
-DTPL_LAPACK_LIBRARIES=/usr/lib/x86_64-linux-gnu/libopenblas.so \
-DTPL_ENABLE_Eigen:BOOL=ON \
-DTPL_Eigen_INCLUDE_DIRS=/usr/include/eigen3 \
-DTrilinos_ENABLE_DataTransferKit=ON \
-DDataTransferKit_ENABLE_DBC=ON \
-DDataTransferKit_ENABLE_TESTS=ON \
-DDataTransferKit_ENABLE_EXAMPLES=OFF \
-DDataTransferKit_ENABLE_ClangFormat=OFF \
-DTPL_ENABLE_BoostLib:BOOL=OFF \
-DBUILD_SHARED_LIBS:BOOL=ON
$ make && sudo make install

```

Again, this assumes root access, adjust this based on your situation. **DTK2** needs to lie in the root directory of **Trilinos** and be turned on through switches `DTrilinos_EXTRA_REPOSITORIES` and `DTrilinos_ENABLE_DataTransferKit`. The environment var `MOAB_ROOT` is where you install **MOAB**.

Note: We recommend that install **DTK2** from our [unifem forked repo](#) or my [personal forked repo](#) since we may add/modify the source codes to make **DTK2** more advanced.

2.3 Install ParPyDTK2

Once you have the dependencies setup, installing ParPyDTK2 can be very easy. The easiest way is through PyPI:

```
$ sudo pip3 install parpydtk2
```

However, this assumes that ParPyDTK2 can find **MOAB** and **DTK2** on the system. Automatically, with different specifications of `install` command, ParPyDTK2 can add different paths in order to locate **MOAB** and **DTK2**.

```
$ pip3 install parpydtk2 --user
```

will assume **MOAB** and **DTK2** can be found in `USER_BASE/{include,lib}`.

```
$ pip3 install parpydtk2 --prefix=...
```

will allow ParPyDTK2 to search **MOAB** and **DTK2** under the `prefix` directory.

The preferred way is to define the environment variables `PARPYDTK2_MOAB_ROOT` and `PARPYDTK2_DTK_ROOT` before you do `pip install`. For instance,

```
$ export PARPYDTK2_MOAB_ROOT=/path/to/moab/root
$ export PARPYDTK2_DTK_ROOT=/path/to/dtk/root
$ pip3 install parpydtk2 --user
```

Warning: We don't mark `mpi4py` as installation dependency, so you need to install it manually before you install ParPyDTK2. `pip3 install mpi4py` is just fine.

Of course, you can install from source, which can be obtained [here](#). Just make sure you have all Python *dependencies* installed.

```
$ git clone -b parallel https://github.com/chiao45/parpydtk2.git
$ cd parpydtk2
$ python3 setup.py install --user
```

2.4 Using our Docker container

You can try the package through our pre-built [Docker container](#). Two driver scripts are provided in order to easily use the container:

1. `parpydtk2_desktop.py`
2. `parpydtk2_jupyter.py`

The former will launch a desktop environment through VNC, while the latter will run the container as a Jupyter server.

SOME DETAILS

3.1 Global IDs/Handles

MOAB uses global IDs in parallel as well as **DTK2**. Therefore, ParPyDTK2 expects the user to provide this information. For most applications, this can be obtained offline.

The global IDs are unique handles of the vertices in a point cloud. For instance, if you want to distribute two triangles, then each of them has local IDs from 0 to 2, but a unique global ID that ranges from 1 to 4.

Note: ParPyDTK2 uses 1-based indexing for global IDs

3.2 Treatment of Empty Partitions

Both **MOAB** and **DTK2** don't support empty partitions, which occur pretty frequently in practice. For instance, couple a serial solver with a parallel solver, or couple a commercial code with an open-sourced one through socket and the incoming data from the commercial code probably doesn't align with the open-sourced side.

To support empty partitions, we duplicate the first node in the master process to the processes that are empty. However, this is not complete yet, because the user is not aware of this, so that the values between the duplicated nodes are not synchronized. Of course, we don't want to let the user to explicitly handle this extra layer of communication. To resolve this, a member function, called `resolve_empty_partitions()`, is added to class `IMeshDB` and must be called **collectively** whenever the user updates the field values.

A DEMO

Here, we show a demo for transferring solutions between two meshes of the unit square. We let the blue mesh participant run in parallel with two cores, while the green side is treated as serial mesh.

```
1 import numpy as np
2 from mpi4py import MPI
3 from parpydtk2 import *
4
5 comm = MPI.COMM_WORLD
6
7 blue = IMeshDB(comm)
8 green = IMeshDB(comm)
9 assert comm.size == 2
10 rank = comm.rank
```

For demo purpose, we assign the meshes globally.

```
12 # create blue meshes on all processes
13 cob = np.empty(shape=(16, 3), dtype=float, order='C')
14 dh = 0.3333333333333333
15 index = 0
16 x = 0.0
17 for i in range(4):
18     y = 0.0
19     for j in range(4):
20         cob[index] = np.asarray([x, y, 0.0])
21         index += 1
22         y += dh
23     x += dh
24
25 # global IDs, one based
26 bgids = np.arange(16, dtype='int32')
27 bgids += 1
```

The blue side has 16 nodes, the following is for the green side:

```
29 # create green meshes on all processes
30 cog = np.empty(shape=(36, 3), dtype=float, order='C')
31 dh = 0.2
32 index = 0
33 x = 0.0
34 for i in range(6):
35     y = 0.0
36     for j in range(6):
37         cog[index] = np.asarray([x, y, 0.0])
```

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```

38         index += 1
39         y += dh
40         x += dh
41
42     # global IDs
43     ggids = np.arange(36, dtype='int32')
44     ggids += 1

```

The green participant has 36 nodes. The next step is to put the data in the two mesh databases.

```

46 # creating the mesh database
47 blue.begin_create()
48 # local vertices and global IDs
49 lcob = cob[8 * rank:8 * (rank + 1), :].copy()
50 lbgsids = bgids[8 * rank:8 * (rank + 1)].copy()
51 blue.create_vertices(lcob)
52 blue.assign_gids(lbgsids)
53 blue.create_field('b')
54
55 # do not use trivial global ID strategy
56 blue.finish_create(False)

```

As we can see, we equally distributed the mesh into the two cores as well as the corresponding global IDs. In line 56, the `False` flag indicates that the mesh database should use the user-provided global IDs.

Warning: Creating vertices and fields, and assigning global IDs must be called within `begin_create()` and `finish_create()`! Otherwise, exceptions are thrown.

Here is the treatment for the “serial” participant:

```

58 # NOTE that green is assumed to be serial mesh
59 green.begin_create()
60 # only create on master rank
61 if not rank:
62     green.create_vertices(cog)
63     green.create_field('g')
64 # since green is serial, we just use the trivial global IDs
65 green.finish_create() # empty partition is resolve here
66
67 assert green.has_empty()

```

As we can see, only the master process has data.

Note: The *duplicated* node is handled inside `finish_create()`

With the two participants ready, we can now create our *Mapper*.

```

69 # create our analytical model, i.e. 10+sin(x)*cos(y)
70 bf = 10.0 + np.sin(lcob[:, 0]) * np.cos(lcob[:, 1])
71 gf = np.sin(cog[:, 0]) * np.cos(cog[:, 1]) + 10.0
72
73 # Construct our mapper
74 mapper = Mapper(blue=blue, green=green)

```

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```

75
76 # using mmls, blue radius 1.0 green radius 0.6
77 mapper.method = MMLS
78 mapper.radius_b = 1.0
79 mapper.radius_g = 0.6
80 mapper.dimension = 2
81
82 # Mapper initialization region
83 mapper.begin_initialization()
84 mapper.register_coupling_fields(bf='b', gf='g', direct=B2G)
85 mapper.register_coupling_fields(bf='b', gf='g', direct=G2B)
86 mapper.end_initialization()

```

Line 69-70 just create an analytic model for error analysis. Line 77-80 are for parameters, for this case, we use MMLS (default) with blue side radius 1.0 and green side radius 0.6 for searching.

The important part is from line 83 to 86. Particularly speaking, the function `register_coupling_fields()`. It takes three parameters, where the first two are string tokens that represents the data fields in blue and green. The `direct` is to indicate the transfer direction, e.g. `B2G` stands for blue to green.

Warning: `register_coupling_fields()` must be called within `begin_initialization()` and `end_initialization()`.

```

88 # NOTE that the following only runs on green master mesh
89 if not rank:
90     green.assign_field('g', gf)
91 # Since green is serial and has empty partition, we must call this to
92 # resolve asynchronous values
93 green.resolve_empty_partitions('g')

```

The above section is to assign values on the green participant. Notice that it is a “serial” mesh, so we only assign values on the master process. But resolve the *duplicated* node is needed, this is done in line 93.

Finally, the solution transfer part is pretty straightforward:

```

94 # solution transfer region
95 mapper.begin_transfer()
96 mapper.transfer_data(bf='b', gf='g', direct=G2B)
97 err_b = (bf - blue.extract_field('b'))/10
98 mapper.transfer_data(bf='b', gf='g', direct=B2G)
99 err_g = (gf - green.extract_field('g'))/10
100 mapper.end_transfer()
101
102 comm.barrier()
103
104 print(rank, 'blue L2-error=%.3e' % (np.linalg.norm(err_b)/np.sqrt(err_b.size)))
105 if rank == 0:
106     print(0, 'green L2-error=%.3e' % (np.linalg.norm(err_g)/np.sqrt(err_g.size)))

```

This code can be obtained [here](#) `parallel2serial.py`.

5.1 ParPyDTK2 Python API

5.1.1 Default

Main module interface of ParPyDTK2

`parpydtk2.B2G`

bool – boolean flag of `True` denotes transferring direction from blue to green

`parpydtk2.G2B`

bool – boolean flag of `False` denotes transferring direction from green to blue

`parpydtk2.MMLS`

int – flag (0) represents using *modified moving least square* method

`parpydtk2.SPLINE`

int – flag (1) represents using *spline interpolation* method

`parpydtk2.N2N`

int – flag (2) represents using *nearest node projection* method

`parpydtk2.WENDLAND2`

int – flag (0) represents using *Wendland 2nd-order* RBF weights

`parpydtk2.WENDLAND4`

int – flag (1) represents using *Wendland 4th-order* RBF weights

`parpydtk2.WENDLAND6`

int – flag (2) represents using *Wendland 6th-order* RBF weights

`parpydtk2.WU2`

int – flag (3) represents using *Wu 2nd-order* RBF weights

`parpydtk2.WU4`

int – flag (4) represents using *Wu 4th-order* RBF weights

`parpydtk2.WU6`

int – flag (5) represents using *Wu 6th-order* RBF weights

`parpydtk2.BUHMANN3`

int – flag (6) represents using *Buhmann 3rd-order* RBF weights

`parpydtk2.get_include()`

Get the abs include path

5.1.2 Error Handling

The error handler module

`parpydtk2.error_handle.ERROR_CODE`

int – set this to nonzero values one exceptions have been raised

Examples

```
>>> import parpydtk2 as dtk
>>> try:
...     # your programs here
... except Exception:
...     dtk.error.ERROR_CODE = 1
...     raise
```

5.1.3 Interface Mesh & Mapper

class `parpydtk2.IMeshDB(comm=None)`

Interface mesh database

ParPyDTK2 utilizes MOAB as the underlying mesh database. MOAB is an array based mesh library that is adapted by DTK2. With array based mesh library, the memory usage and computational cost are lower than typical pointer based data structure. The mesh concept in this work is simple since only meshless methods are utilized, the only additional attribute one needs is the *global IDs/handles*, which are used by both MOAB and DTK2. For most applications, the global IDs can be computed offline.

One thing is not directly supported by IMeshDB is I/O. However, since this is a Python module and only points clouds are needed, one can easily use a tool (e.g. [meshio](#)) to load the mesh.

comm

MPI.Comm – MPI communicator

ranks

int – size of comm

rank

int – rank of comm

size

int – point cloud size, i.e. number of vertices

bbox

np.ndarray – local bounding box array of shape (2,3)

gbbox

np.ndarray – global bounding box array of shape (2,3)

Constructor

Parameters **comm** (*MPI.Comm* (optional)) – if no communicator or None is passed in, then `MPI_COMM_WORLD` will be used

Examples

```
>>> # implicit communciator
>>> import parpydtk2 as dtk
>>> mdb = dtk.IMeshDB()
```

```
>>> # explicit communicator
>>> from mpi4py import MPI
>>> import parpydtk2 as dtk
>>> mdb = dtk.IMeshDB(MPI.COMM_WORLD)
```

assign_field(*self*, *unicode field_name*, *ndarray values*)
Assign values to a field

Note: *values* size must be at least *size*dim*

Parameters

- **field_name** (*str*) – name of the field
- **values** (*np.ndarray*) – input source values

See also:

[*extract_field\(\)*](#) extract value from a field

[*size*](#) check the size of a mesh set

assign_gids(*self*, *__Pyx_memviewslice gids*)
Assign global IDs

Internally, both DTK and MOAB use so-called global IDs/handles communications. Each node has its own local IDs/handles and a unique global ID.

Parameters **gids** (*np.ndarray*) – global IDs

See also:

[*create_vertices\(\)*](#) create vertices

[*extract_gids\(\)*](#) extract global IDs

bbox

np.ndarray – local bounding box

The bounding box is stored simply in a 2x3 array, where the first row stores the maximum bounds while minimum bounds for the second row.

Warning: Bounding box is valid only after [*finish_create\(\)*](#).

See also:

[*gbbox*](#) global bounding box

begin_create (*self*)

Begin to create/manupilate the mesh

This function must be called in order to let the mesh database be aware that you will create meshes.

See also:

[`finish_create\(\)`](#) finish creating mesh

comm

MPI.Comm – communicator

create_field (*self*, *unicode field_name*, *int dim=1*)

Create a data field for solution transfer

This is the core function to register a field so that you can then transfer its values to other domains. The `dim` parameter determines the data type of the field. By default, it's 1, i.e. scalar fields. For each node, a tensor of (1x“dim“) can be registered. For instance, to transfer forces and displacements in FSI applications, `dim` is 3 (for 3D problems).

Parameters

- **field_name** (*str*) – name of the field
- **dim** (*int*) – dimension of the field, i.e. scalar, vector, tensor

Examples

```
>>> from parpydtk2 import *
>>> mdb1 = IMeshDB()
>>> mdb1.begin_create()
>>> mdb1.create_field('heat flux')
```

create_vertices (*self*, *__Pyx_memviewslice coords*)

Create a set of coordinates

Note: The `coords` must be C-ordering with `ndim=2!`

Parameters **coords** (*np.ndarray*) – nx3 coordinates in double precision

See also:

[`assign_gids\(\)`](#) assign global IDs

[`extract_vertices\(\)`](#) extract vertex coordinates

Examples

```
>>> from parpydtk2 import *
>>> import numpy as np
>>> mdb1 = IMeshDB()
>>> mdb1.begin_create()
>>> verts = np.zeros((2,3)) # two nodes
```

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```
>>> verts[1][0] = 1.0
>>> mdb1.create_vertices(verts)
```

empty (*self*)

Check if this is an empty partition

extract_field (*self*, *unicode field_name*, *__Pyx_memviewslice buffer=None*, *reshape=False*)

Extract the values from a field

Warning: if *buffer* is passed in, it must be 1D

Parameters

- **field_name** (*str*) – name of the field
- **buffer** (*np.ndarray*) – 1D buffer
- **reshape** (*bool*) – *True* if we reshape the output, only for vectors/tensors

Returns field data values**Return type** *np.ndarray***extract_gids** (*self*)

Extract global IDs/handles

Warning: This function should be called once you have finished *assign_gids()*.

Returns array of *size* that stores the integer IDs**Return type** *np.ndarray***extract_vertices** (*self*)

Extract coordinate

Warning: This function should be called once you have finished *create_vertices()*.

Returns (*nx3*) array that stores the coordinate values**Return type** *np.ndarray***See also:***size* get the mesh size**field_dim** (*self*, *unicode field_name*)

Check the field data dimension

Parameters **field_name** (*str*) – name of the field**Returns** data field dimension of *field_name***Return type** *int*

finish_create (*self*, *trivial_gid=True*)
 finish mesh creation

This method finalizes the interface mesh database by communicating the bounding boxes and empty partitions. Also, setting up the DTK managers happens here.

Warning: You must call this function once you have done with manipulating the mesh, i.e. vertices and global IDs.

Parameters **trivial_gid** (*bool*) – *True* if we use MOAB trivial global ID computation

Notes

By `trivial_gid`, it means simply assigning the global IDs based on the size of the mesh. This is useful in serial settings or transferring solutions from a serial solver to a partitioned one.

gbbox

np.ndarray – global bounding box

The bounding box is stored simply in a 2x3 array, where the first row stores the maximum bounds while minimum bounds for the second row.

Warning: Bounding box is valid only after `finish_create()`.

See also:

[*bbox*](#) local bounding box

has_empty (*self*)

Check if an empty partition exists

has_field (*self*, *unicode field_name*)

Check if a field exists

Parameters **field_name** (*str*) – name of the field

Returns *True* if this meshdb has *field_name*

Return type *bool*

rank

int – get the rank

ranks

int – Get the communicator size

resolve_empty_partitions (*self*, *unicode field_name*)

Resolve asynchronous values on empty partitions

ParPyDTK2 doesn't expect `assign_field()` should be called collectively. Therefore, a collective call must be made for resolving empty partitions.

Warning: This must be called collectively even on empty partitions

Note: You should call this function following assignment

Parameters `field_name` (*str*) – name of the field

Examples

```
>>> if rank == 0:
...     mdb.assign_field('flux', values)
>>> mdb.resolve_empty_partitions('flux')
```

Notes

This API is not available in C++ level, therefore, one needs to implement this if he/she wants to use the C++ API.

size

int – Get the size of a set

class `parpydtk2.Mapper` (*blue, green, profiling=True*)

DTK2 wrapper

The meshless methods in DTK2, including *modified moving least square*, *spline interpolation* and *nearest node projection* methods are wrapped within this class. The most advanced method is the MMLS fitting, which is my personal recommendaion.

blue_mesh

IMeshDB – blue mesh participant

green_mesh

IMeshDB – green mesh participant

comm

MPI.Comm – MPI communicator

ranks

int – size of comm

rank

int – rank of comm

dimension

int – spatial dimension

method

int – method flag, either MMLS, SPLINE, or N2N

basis

int – flag of basis function for weighting schemes used by MMLS and SPLINE

knn_b

int – k-nearest neighborhood for searching on blue_mesh

knn_g

int – k-nearest neighborhood for searching on green_mesh

radius_b*float* – radius used for searching on blue_mesh**radius_g***float* – radius used for searching on green_mesh

Constructor

Parameters

- **blue_mesh** (*IMeshDB*) – blue mesh participant
- **green_mesh** (*IMeshDB*) – green mesh participant
- **profiling** (*bool (optional)*) – whether or not do timing report, default is `True`.

Examples

```
>>> from parpydtk2 import *
>>> blue, green = IMeshDB(), IMeshDB()
>>> # initialize blue and green
>>> mapper = Mapper(blue=blue, green=green)
>>> # do work with mapper
```

basis*int* – Get the basis function flag**See also:***method* get the method tag**begin_initialization** (*self*)

Initialization starter

This is a must-call function in order to indicate mapper that you are about to initialize/register coupling fields

See also:*register_coupling_fields()* register coupled fields*end_initialization()* finish initialization**begin_transfer** (*self*)

Transfer starter

This is a must-call function to inidate the beginning of a transferring block

See also:*end_transfer()* transfer closer*tranfer_data()* transfer a coupled data fields**blue_mesh***IMeshDB* – blue mesh**comm***MPI.Comm* – Get the communicator**See also:**

ranks get the total communicator size

rank “my” rank

dimension

int – Get the problem dimension

Note: this is the spacial dimension

end_initialization (*self*)

Initialization closer

This is a must-call function in order to tell the mapper we are ready

See also:

begin_initialization() initialization starter

end_transfer (*self*)

Transfer closer

This is a must-call function to indicate we have finished a sequence of transferring requests

See also:

begin_transfer() transfer starter

green_mesh

IMeshDB – green mesh

has_coupling_fields (*self, unicode bf, unicode gf, bool direct*)

Check if a coupled fields exists

Returns True if (bf,gf) exists in the `direct` direction

Return type bool

knn_b

int – KNN of blue mesh

Note: if blue does not use KNN, then -1 returned

See also:

knn_g green knn

knn_g

int – KNN of green mesh

Note: if green does not use KNN, then -1 returned

See also:

knn_b blue knn

method*int* – Get the method tag**See also:***basis* the basis function and order attribute**radius_b***float* – physical domain radius support for blue mesh

Note: if blue does not use RBF-search, then -1.0 returned

See also:*radius_g* green radius**radius_g***float* – physical domain radius support for green mesh

Note: if green does not use RBF-search, then -1.0 returned

See also:*radius_b* blue radius**rank***int* – Check “my” rank**See also:***ranks* get the total communicator size*comm* MPI communicator**ranks***int* – Check the total process number**See also:***rank* “my” rank**register_coupling_fields** (*self, unicode bf, unicode gf, bool direct*)

register a coupled fields

Note: we use boolean to indicate direction

Parameters

- **bf** (*str*) – blue mesh field name
- **gf** (*str*) – green mesh field name
- **direct** (*bool*) – *True* for blue->green, *False* for the opposite

See also:

`transfer_data()` transfer a coupled data fields

`set_matching_flag_n2n(self, bool matching)`
Set the matching flag for N2N

Note: this function will not throw even if you dont use n2n

Parameters **matching** (*bool*) – *True* if the interfaces are matching

`transfer_data(self, unicode bf, unicode gf, bool direct)`
Transfer (bf, gf) in the `direct` direction

Parameters

- **bf** (*str*) – blue mesh field name
- **gf** (*str*) – green mesh field name
- **direct** (*bool*) – *True* for blue->green, *False* for the opposite

See also:

`register_coupling_fields()` register coupled fields

5.2 ParPyDTK2 C++ API

5.2.1 Common Definitions

group **common**

Defines

handle_moab_error (__ret)
macro to handle moab error

throw_error (__msg)
throw `runtime_error` exception

throw_error_if (__cond, __msg)
conditionally error throw

throw_noimpl (__what)
throw not implemented feature error

throw_noimpl_if (__cond, __what)
throw not implemented feature error with condition

show_warning (__msg)
log warning message in `stderr`

show_warning_if (__cond, __msg)
log warning with condition

show_experimental (__msg)
log experimental warning in stderr

show_experimental_if (__cond, __msg)
log experimental warning in stderr with condition

show_info (__msg, __rank)
show information in parallel

show_info_master (__msg, __rank)
show information only on master rank

streamer (__rank)
streaming message with specific rank

streamer_master (__rank)
streaming messages only on the master process

Typedefs

typedef entity_t
MOAB entity handle.

Enums

enum [anonymous]
root set

Values:

root_set = 0

5.2.2 Field Variables

class FieldData
a representation of MOAB tag for field data

Public Functions

FieldData (moab::Core &mdb, **const** std::string &field_name, int dim = 1)
constructor with moab instance

Note *set* is not the same as mesh set in MOAB

Parameters

- mdb: moab instance
- field_name: field name
- dim: field dimension

void **assign** (**const** moab::Range &range, **const** double *values)
assign values

Parameters

- `range`: entity ranges, for this work, it should be vertices
- `values`: data values, for vector/tensor, C order is expected

void **assign_1st** (**const** moab::Range &*range*, **const** double **values*)
 assign to first node

This function is used by Python for handling empty partitions

Parameters

- `range`: entity ranges
- `values`: values for the first node, at least size of dim

void **extract** (**const** moab::Range &*range*, double **values*) **const**
 extract values

Parameters

- `range`: entity ranges, for this work, it should be vertices
- `values`: data values, for vector/tensor, C order is expected

void **extract_1st** (**const** moab::Range &*range*, double **values*) **const**
 extract the value from the first node

Parameters

- `range`: entity ranges
- `values`: values for the first node, at least size of dim

operator const std::string& () const
 brief implicitly cast to string

int **dim () const**
 check the dimension

int **set () const**
 check the set ID

const moab::Tag &**tag () const**
 get MOAB tag

Protected Attributes

moab::Core &**mdb_**
 reference to moab instance

std::string **fn_**
 field name

int **set_**
 set count

```

int dim_
    field dimension

moab::Tag tag_
    moab tag

class FieldDataSet
    a set of field data

```

Public Types

```

typedef base_t::iterator iterator
    iterator type

typedef base_t::const_iterator const_iterator
    constant iterator

```

Public Functions

```

virtual ~FieldDataSet ()
    destructor

bool has_field (const std::string &fn) const
    check if a field exist

```

Parameters

- *fn*: field name

```

void create (moab::Core &mdb, const std::string &field_name, int dim = 1)
    create an data field

```

Note *set* is not the same as mesh set in MOAB

Parameters

- *mdb*: moab data base
- *field_name*: field name
- *dim*: field dimension

```

FieldData &operator [] (const std::string &fn)
    get a reference to a field data

```

Note this overloads the base operator[]

Parameters

- *fn*: field name

```

const FieldData &operator [] (const std::string &fn) const
    get a const reference to a field data

```

Parameters

- *fn*: field name

```

iterator begin ()
    get the first iterator

iterator end ()
    get the end iterator

const_iterator begin () const
    get the constant iterator

const_iterator end () const
    get the constant end iterator

const_iterator cbegin () const
    get the constant iterator

const_iterator cend () const
    get the constant end iterator

```

Protected Attributes

```

base_t fs_
    fields

```

Private Types

```

typedef std::unordered_map<std::string, FieldData *> base_t
    data structure

```

5.2.3 Interface Mesh Database

```

class IMeshDB
    interface mesh database, build on top of MOAB

```

imesh_py_interface

```

void begin_create ()
    begin to create mesh

void create_vset ()
    create a new vertex set

void create_vertices (int nv, const double *coords)
    create vertices

```

Parameters

- *nv*: number of vertices
- *coords*: coords values

```

void extract_vertices (double *coords) const
    extract assigned coordinates

```

Note coords must be at least $n*3$ where n is the size of the mesh

Parameters

- `coords`: coordinates

void **assign_gids** (int *nv*, **const** int **gids*)
assign global IDs

Note *gids* should be one-based indices

Parameters

- *nv*: number of local vertices
- *gids*: global IDs

void **extract_gids** (int **gids*) **const**
extract global IDs

Parameters

- *gids*: global IDs

void **finish_create** (bool *trivial_gid* = true)
finish manipulating the mesh

NOTE that if your input mesh is element-based partition and the vertices you create are mesh nodes, then you have to specify the correct global IDs in parallel. However, if the coordinates are face centres, then the global IDs can be trivially computed by MOAB since there are no shared entities cross different processes.

Parameters

- *trivial_gid*: *true* if we let MOAB to compute the GID

bool **empty** () **const**
check if empty partition

bool **has_empty** () **const**
check if any of the process has an empty partition

const std::vector<int> &**m2s** () **const**
get a reference to the m2s pattern

Note This is used in Python level as “private” thus having “_”

int **size** () **const**
check mesh size

void **get_bbox** (double **v*) **const**
get bounding box

Parameters

- *v*: values

void **get_gbbox** (double **v*) **const**
get global bounding box

Parameters

- *v*: values

void **create_field**(**const** std::string &*field_name*, int *dim* = 1)
create a field

Parameters

- *field_name*: field name
- *dim*: field dimension

bool **has_field**(**const** std::string &*field_name*) **const**
check if we have a field

Parameters

- *field_name*: field name

int **field_dim**(**const** std::string &*field_name*) **const**
check field dimension

Parameters

- *field_name*: field name

void **assign_field**(**const** std::string &*field_name*, **const** double **values*)
assign a value to a field

Parameters

- *field_name*: field name
- *values*: field data values

void **_assign_1st**(**const** std::string &*field_name*, **const** double **values*)
assign to the first node

This function is used by Python to resolve the issues when empty empty partitions happen. Therefore, this function has an “_” prefix to indicate “private” usage!

Parameters

- *field_name*: field name
- *values*: field data values, at least size of field dimension

void **extract_field**(**const** std::string &*field_name*, double **values*) **const**
extract value

Parameters

- *field_name*: field name
- *values*: field data values

```
void _extract_1st (const std::string &field_name, double *values) const
    extract first value
```

This function is used by Python to resolve the issues when empty empty partitions happen. Therefore, this function has an “_” prefix to indicate “private” usage!

Parameters

- field_name: field name
- values: field data values

Public Functions

```
IMeshDB (MPI_Comm comm = MPI_COMM_WORLD)
    constructor with communicator
```

Parameters

- comm: communicator

```
int ranks () const
    get total ranks
```

```
int rank () const
    get my rank
```

```
MPI_Comm comm () const
    get the communicator
```

```
Teuchos::RCP<moab::ParallelComm> pcomm () const
    get mesh
```

```
std::vector<DataTransferKit::MoabManager> &mangers ()
    get the manger
```

```
void set_dimension (int dim)
    set geometry dimension
```

Parameters

- dim: dimension

```
bool ready () const
    check if ready
```

```
dtk_field_t &dtk_fields ()
    get the dtk fields
```

Protected Types

```
typedef std::unordered_map<std::string, std::pair<int, Teuchos::RCP<Tpetra::MultiVector<double, int, DataTransferKit::Supp
    handy typedef
```

Protected Attributes

moab::Core **mdb_**
moab instance

Teuchos::RCP<moab::ParallelComm> **par_**
moab parallel interface

std::vector<*entity_t*> **vsets_**
vertex sets

std::vector<moab::Range> **locals_**
local vertex range

std::vector<std::array<double, 6>> **bboxes_**
bounding boxes

std::vector<std::array<double, 6>> **gbboxes_**
global bounding boxes

FieldDataSet **fields_**
field data set

bool **created_**
flag to indicate whether users are done with creating mesh

moab::Tag **gidtag_**
global ID tag

moab::Tag **parttag_**
partition tag

bool **usergid_**
flag to indicate if we have user computed global ID

std::vector<DataTransferKit::MoabManager> **mngrs_**
DTK MOAB manager.

dtk_field_t **dtkfields_**
DTK multi vector for MOAB tags.

bool **empty_**
check empty partition

bool **has_empty_**
check if any process is empty

std::vector<int> **m2s_**
comm pattern for master2slaves for handling empty partitions

Private Functions

void **init_** (bool *del* = false)
helper for clean up mesh

Parameters

- *del*: whether or not delete mesh

void **reset_vecs_** ()
handle all vectors

void **init_bbox_** (int *i* = -1)
initialize empty bounding boxes

Parameters

- *i*: index if < 0 then init all

void **cmpt_bboxes_** ()
compute all bounding box

5.2.4 Solution Mapper

class Mapper

the mapper interface for interface solution transfer

mapper_py_interface

int **ranks** () **const**
get the ranks

int **rank** () **const**
get my rank

MPI_Comm **comm** () **const**
get the communicator

void **set_dimension** (int *dim*)
set dimension

Parameters

- *dim*: geometry dimension

void **use_mmls** ()
use moving least square, this is the default method

See *use_spline, use_n2n*

void **use_spline** ()
use spline interpolation method

See *use_mmls, use_n2n*

void **use_n2n** (bool *matching* = false)
use node 2 node project

See *use_mmls, use_spline*

Parameters

- *matching*: are the interfaces mathing?

void **set_basis** (int *basis*)
set basis function, default is Wendland 4th order

See BasisFunctions

Parameters

- *basis*: basis function and order

void **use_knn_b** (int *knn*)
use knn for blue mesh

Parameters

- *knn*: number of nearest neighbors

void **use_knn_g** (int *knn*)
use knn for green mesh

Parameters

- *knn*: number of nearest neighbors

void **use_radius_b** (double *r*)
use radius for blue

See use_knn

Parameters

- *r*: physical domain radius support

void **use_radius_g** (double *r*)
use radius for green

See use_knn

Parameters

- *r*: physical domain radius support

int **check_method** () **const**
check method

int **check_basis** () **const**
check basis

int **knn_b** () **const**
check blue knn

Note if blue does not use knn, then negative value returned

int **knn_g** () **const**
check green knn

double **radius_b** () **const**
check blue radius

Note if blue does not use radius, then -1.0 is returned

```
double radius_g() const
    check green radius
```

```
int dimension() const
    get the dimension
```

```
std::shared_ptr<IMeshDB> blue_mesh() const
    get blue mesh
```

```
std::shared_ptr<IMeshDB> green_mesh() const
    get green mesh
```

```
void begin_initialization()
    begin initialization
```

```
void register_coupling_fields(const std::string &bf, const std::string &gf, bool direct)
    register coupling fields
```

Parameters

- *bf*: blue meshdb field data
- *gf*: green meshdb field data
- *direct*: *true* for b->g, *false* for g->b

```
bool has_coupling_fields(const std::string &bf, const std::string &gf, bool direct)
    check if a coupling data fields exists
```

Parameters

- *bf*: blue meshdb field data
- *gf*: green meshdb field data
- *direct*: *true* for b->g, *false* for g->b

```
void end_initialization()
    end initialization
```

```
void begin_transfer()
    begin to transfer data
```

```
void transfer_data(const std::string &bf, const std::string &gf, bool direct)
    transfer data
```

Parameters

- *bf*: blue meshdb field data
- *gf*: green meshdb field data
- *direct*: *true* for b->g, *false* for g->b

```
void end_transfer()
    end transfer
```

Public Functions

Mapper (std::shared_ptr<*IMeshDB*> *B*, std::shared_ptr<*IMeshDB*> *G*, **const** std::string &*version* = "",
const std::string &*date* = "", bool *profiling* = true)
 constructor

Parameters

- *B*: input blue mesh
- *G*: input green mesh
- *version*: passed from Python interface
- *date*: passed from Python interface
- *profiling*: whether do simple profiling, i.e. wtime

Protected Attributes

std::shared_ptr<*IMeshDB*> **B_**
 blue mesh

std::shared_ptr<*IMeshDB*> **G_**
 green mesh

int **dim_**
 dimension

bool **ready_**
 flag to indicate the mapper is ready for transferring

bool **profiling_**
 whether do simple profiling

double **timer_**
 a simple timer buffer

std::unique_ptr<Teuchos::ParameterList> **opts_[2]**
 parameter list

std::map<std::pair<std::string, std::string>, Teuchos::RCP<DataTransferKit::MapOperator>> **operators_[2]**
 transfer operators

Protected Static Attributes

DataTransferKit::MapOperatorFactory **factory_**
 map factory

Private Functions

void **init_parlist_** ()
 initialize parameter list

template <bool *_Dir*, typename *_V*>
 void **set_search** (**const** std::string &*type*, **const** std::string &*value*, **const** *_V* &*v*)
 helper for set local search

Template Parameters

- `_Dir`: direction
- `_V`: value type

Parameters

- `type`: search type
- `value`: tag for value
- `v`: actual value

```
template <bool _Dir, typename _V>  
_V get_search (const std::string &type, const std::string &value, const _V &dft) const  
    helper for get search info
```

Template Parameters

- `_Dir`: direction
- `_V`: value type

Parameters

- `type`: search type
- `value`: tag for value
- `dft`: default value

Private Static Functions

```
static std::string parse_list_ (Teuchos::ParameterList &list)  
    parse and formatting a parameter list
```

Parameters

- `list`: parameter list

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