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Zoltan: Parallel Partitioning, Load Balancing and Data-Management Services

User's Guide

The Zoltan Team

Sandia National Laboratories

[Erik Boman](#)

Cedric Chevalier

[Karen Devine](#)

Vitus Leung

Sivasankaran Rajamanickam

Lee Ann Riesen

Michael Wolf

Ohio State University

[Umit Catalyurek](#)

[Doruk Bozdag](#)

Past Zoltan Contributors

Sandia National Laboratories:

Robert Heaphy

[Bruce Hendrickson](#)

Matthew St. John

Courtenay Vaughan

Williams College

[James Teresco](#)

National Institute of Standards and Technology

[William F. Mitchell](#)

Rensselaer Polytechnic Institute

Jamal Faik

Luis Gervasio

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Introduction

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Project Motivation

Over the past decade, parallel computers have been used with great success in many scientific simulations. While differing in their numerical methods and details of implementation, most applications successfully parallelized to date are "static" applications. Their data structures and memory usage do not change during the course of the computation. Their inter-processor communication patterns are predictable and non-varying. And their processor workloads are predictable and roughly constant throughout the simulation. Traditional finite difference and finite element methods are examples of widely used static applications.

However, increasing use of "dynamic" simulation techniques is creating new challenges for developers of parallel software. For example, adaptive finite element methods refine localized regions the mesh and/or adjust the order of the approximation on individual elements to obtain a desired accuracy in the numerical solution. As a result, memory must be allocated dynamically to allow creation of new elements or degrees of freedom. Communication patterns can vary as refinement creates new element neighbors. And localized refinement can cause severe processor load imbalance as elemental and processor work loads change throughout a simulation.

Particle simulations and crash simulations are other examples of dynamic applications. In particle simulations, scalable parallel performance depends upon a good assignment of particles to processors; grouping physically close particles within a single processor reduces inter-processor communication. Similarly, in crash simulations, assignment of physically close surfaces to a single processor enables efficient parallel contact search. In both cases, data structures and communication patterns change as particles and surfaces move. Re-partitioning of the particles or surfaces is needed to maintain geometric locality of objects within processors.

We developed the Zoltan library to simplify many of the difficulties arising in dynamic applications. Zoltan is a collection of data management services for unstructured, adaptive and dynamic applications. It includes a suite of parallel partitioning algorithms, data migration tools, parallel graph coloring tools, distributed data directories, unstructured communication services, and dynamic memory management tools. Zoltan's data-structure neutral design allows it to be used by a variety of applications without imposing restrictions on application data structures. Its object-based interface provides a simple and inexpensive way for application developers to use the library and researchers to make new capabilities available under a common interface.

The Zoltan Toolkit

The Zoltan Library contains a number of tools that simplify the development and improve the performance of parallel, unstructured and adaptive applications. The library is organized as a toolkit, so that application developers can use as little or as much of the library as desired. The major packages in Zoltan are listed below.

- A suite of [dynamic load balancing and parallel repartitioning](#) algorithms, including geometric, hypergraph and graph methods; switching between algorithms is easy, allowing straightforward comparisons of algorithms in

applications.

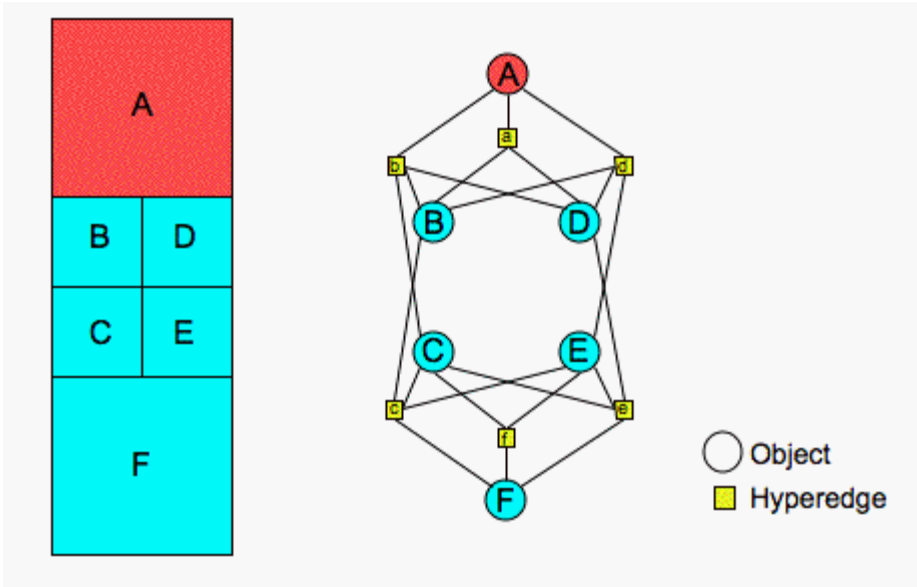
- [Data migration tools](#) for moving data from old partitions to new one.
- [Parallel graph coloring tools](#) with both distance-1 and distance-2 coloring.
- [Distributed data directories](#): scalable (in memory and computation) algorithms for locating needed off-processor data.
- An [unstructured communication package](#) that insulates users from the details of message sends and receives.
- [Dynamic memory management tools](#) that simplify dynamic memory debugging on state-of-the-art parallel computers.
- A sample application [zdrive](#). It allows algorithm developers to test changes to Zoltan without having to run Zoltan in a large application code. Application developers can use the *zdrive* code to see examples of function calls to Zoltan and the implementation of query functions.

Terminology

Our design of Zoltan does not restrict it to any particular type of application. Rather, Zoltan operates on uniquely identifiable data items that we call *objects*. For example, in finite element applications, objects might be elements or nodes of the mesh. In particle applications, objects might be particles. In linear solvers, objects might be matrix rows or non-zeros.

Each object must have a unique *global identifier (ID)* represented as an array of unsigned integers. Common choices include global numbers of elements (nodes, particles, rows, and so on) that already exist in many applications, or a structure consisting of an owning processor number and the object's local-memory index. Objects might also have local (to a processor) IDs that do not have to be unique globally. Local IDs such as addresses or local-array indices of objects can improve the performance (and convenience) of Zoltan's interface to applications.

We use a simple example to illustrate the above terminology. On the left side of the figure [below](#), a simple finite element mesh is presented.



The blue and red shading indicates the mesh is partitioned for two processors. An application must provide information about the current mesh and partition to Zoltan. If, for example, the application wants Zoltan to perform operations on the elements of the mesh, it must provide information about the elements when Zoltan asks for object information.

In this example, the elements have unique IDs assigned to them, as shown by the letters in the elements. These unique letters can be used as global IDs in Zoltan. In addition, on each processor, local numbering information may be available. For instance, the elements owned by a processor may be stored in arrays in the processor's memory. An

element's local array index may be provided to Zoltan as a local ID.

For geometric algorithms, the application must provide coordinate information to Zoltan. In this example, the coordinates of the mid-point of an element are used.

For hypergraph- and graph-based algorithms, information about the connectivity of the objects must be provided to Zoltan. In this example, the application may consider elements connected if they share a face. A hypergraph representing this problem is then shown to the right of the mesh. A *hyperedge* exists for each object (squares labeled with lower-case letters corresponding to the related object). Each hyperedge connects the object and all of its face neighbors. The hyperedges are passed to Zoltan with a label (in this example, a lower-case letter) and a list of the object IDs in that hyperedge.

Graph connections, or *edges*, across element faces may also specified. Connectivity information is passed to Zoltan by specifying a neighbor list for an object. The neighbor list consists of the global IDs of neighboring objects and the processor(s) currently owning those objects. Because relationships across faces are bidirectional, the graph edge lists and hypergraph hyperedge lists are nearly identical. If, however, information flowed to, say, only the north and east edge neighbors of an element, the hypergraph model would be needed, as the graph model can represent only bidirectional relationships. In this case, the hyperedge contents would include only the north and east neighbors; they would exclude south and west neighbors.

The table below summarizes the information provided to Zoltan by an application for this finite element mesh. Information about the objects includes their global and local IDs, geometry data, hypergraph data, and graph data.

	Object IDs		Geometry Data	Graph Data	
Processor	Global	Local	(coordinates)	Neighbor Global ID List	Neighbor Processor List
Red	A	0	(2,10)	B,D	Blue
Blue	B	0	(1,7)	A,C,D	Red,Blue,Blue
	C	1	(1,5)	B,E,F	Blue,Blue,Blue
	D	2	(3,7)	A,B,E	Red,Blue,Blue
	E	3	(3,5)	C,D,F	Blue,Blue,Blue
	F	4	(2,2)	C,E	Blue,Blue

Hyperedge Data	
Hyperedge ID	Hyperedge contents
a	A,B,D
b	A,B,C,D
c	B,C,E,F
d	A,B,D,E
e	C,D,E,F
f	C,E,F

Zoltan Design

To make Zoltan easy to use, we do not impose any particular data structure on an application, nor do we require an application to build a particular data structure for Zoltan. Instead, Zoltan uses a [callback function interface](#), in which Zoltan queries the application for needed data. The application must provide simple functions that answer these queries.

To keep the application interface simple, we use a small set of [callback functions](#) and make them easy to write by requesting only information that is easily accessible to applications. For example, the most basic partitioning algorithms require only four callback functions. These functions return the number of objects owned by a processor, a list of weights and IDs for owned objects, the problem's dimensionality, and a given object's coordinates. More sophisticated graph-based partitioning algorithms require only two additional callback functions, which return the number of edges per object and edge lists for objects.

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Using the Zoltan library

This section contains information needed to use the Zoltan library with applications:

- [System requirements](#)
- [Building the Zoltan library](#)
- [Testing the Zoltan library](#)
- [Reporting bugs](#) in Zoltan
- [Incorporating Zoltan into Applications](#)
- [Building applications](#) that use Zoltan
- [Data types](#) for global and local IDs
- [C++ interface](#)
- [F90 interface](#)

System Requirements

Zoltan was designed to run on parallel computers and clusters of workstations. The most common builds and installations of Zoltan use the following:

- ANSI C or C++ compiler.
- [MPI](#) library for message passing (version 1.1 or higher), such as MPICH, OpenMPI or LAM.
- A Unix-like operating system (e.g., Linux or Mac OS X) and *gmake* (GNU Make) are recommended to build the library.
- A Fortran90 compatible compiler is required if you wish to [use Zoltan with Fortran applications](#).

Zoltan has been tested on a variety of platforms, including Linux, Mac OS X, a variety of clusters, and Sandia's ASC RedStorm machine. Builds for Windows platforms are available as part of the [Trilinos CMake build system](#).

Building the Zoltan Library

The Zoltan library is implemented in ANSI C and can be compiled with any ANSI C compiler. In Zoltan, there are two build environments currently supported: an [Autotools build environment](#) and a [CMake build environment](#) used with the [Trilinos](#) framework. The Autotools build environment can be used to build Zoltan in a stand-alone installation; the CMake build environment must be used within Trilinos.

Using Autotools to Build Zoltan

Users should not run autotools directly in the main Zoltan directory; rather they should create a build-directory (e.g., a subdirectory of the main Zoltan directory) in which they configure and build Zoltan. Say, for example, a user creates a directory called BUILD_DIR in the Zoltan directory. Then, to configure and build zoltan, the user would

```
cd zoltan/BUILD_DIR
../configure {options described below}
make everything
make install
```

Options to the configure command allow paths to third-party libraries such as ParMETIS, PT-Scotch and PaToH to be specified. Building with MPI compilers (e.g., mpicc) is the default for Autotools builds of Zoltan; many options allow specification of particular MPI paths and compilers.

Users desiring a [Fortran90 interface](#) to Zoltan must specify the "--enable-f90interface" option.

All options can be seen with the following command issued in the zoltan/BUILD_DIR directory:

```
../configure --help
```

The following script is an example of configuration and build commands using Autotools. It specifies that Zoltan should be built with both the [ParMETIS](#) and [PT-Scotch](#) interfaces. Paths to both ParMETIS and PT-Scotch are given. The prefix option states where Zoltan should be installed; in this example, Zoltan's include files will be installed in /homes/username/zoltan/BUILD_DIR/include, and the libraries will be installed in /homes/username/zoltan/BUILD_DIR/lib. This examples assumes the script is run from /homes/username/zoltan/BUILD_DIR.

```
#
../configure \
--prefix=/homes/username/zoltan/BUILD_DIR \
--with-gnumake \
--with-scotch \
--with-scotch-incdir="/Net/local/proj/zoltan/arch/all/src/Scotch5" \
--with-scotch-libdir="/Net/local/proj/zoltan/arch/linux64/lib/openmpi/Scotch5" \
--with-parmetis \
--with-parmetis-incdir="/Net/local/proj/zoltan/arch/all/src/ParMETIS3" \
--with-parmetis-libdir="/Net/local/proj/zoltan/arch/linux64/lib/openmpi/ParMETIS3"
make everything
make install
```

More examples are in the directory zoltan/SampleConfigurationScripts.

After the configuration is done in the build directory, object files and executables can be removed with *make clean*; the same configuration will be used for subsequent builds. Configuration information is removed with *make distclean*.

Using CMake to Build Zoltan

Zoltan can be built as part of the Trilinos framework using the CMake build system. CMake builds will succeed only when Zoltan is in the Trilinos directory structure (as when downloaded with Trilinos). Users should not run CMake directly in the main Zoltan directory; rather they should create a build-directory (e.g., a subdirectory of the main Trilinos directory) in which they configure and build Zoltan. Say, for example, a user creates a directory called BUILD_DIR in the Trilinos directory. Then, to configure and build zoltan, the user would

```
cd Trilinos/BUILD_DIR
cmake \
-D Trilinos_ENABLE_ALL_PACKAGES:BOOL=OFF \
-D Trilinos_ENABLE_Zoltan:BOOL=ON \
{options described below} \
..
make
make install
```

Serial builds are the default in Trilinos; for serial builds, Zoltan builds and links with the siMPI library provided by Pat Miller in the Zoltan distribution. More commonly, Zoltan users desire **parallel** builds with MPI libraries such as OpenMPI or MPICH. For such builds, users must specify CMake option

```
-D TPL_ENABLE_MPI:BOOL=ON
```

Trilinos also defaults to using a Fortran compiler, but Fortran is not required to build Zoltan; the option to disable this check is

```
-D Trilinos_ENABLE_Fortran:BOOL=OFF
```

Other options to the cmake command allow paths to third-party libraries such as ParMETIS, PT-Scotch and PaToH to be specified.

Users desiring a [Fortran90 interface](#) to Zoltan must specify the option

```
-D Zoltan_ENABLE_F90INTERFACE:BOOL=ON
```

All options can be seen with the following command issued in the Trilinos/BUILD_DIR directory:

```
rm CMakeCache.txt
cmake -LAH -D Trilinos_ENABLE_Zoltan:BOOL=ON ..
```

The following script is an example of configuration and build commands using CMake. It specifies that Zoltan should be built with both the [ParMETIS](#) and [PT-Scotch](#) interfaces. Paths to both ParMETIS and PT-Scotch are given. The prefix option states where Zoltan should be installed; in this example, Zoltan's include files will be installed in /homes/username/Trilinos/BUILD_DIR/include, and the libraries will be installed in /homes/username/Trilinos/BUILD_DIR/lib. This examples assumes the script is run from /homes/username/Trilinos/BUILD_DIR.

```
#
cmake \
-D CMAKE_INSTALL_PREFIX:FILEPATH="/home/username/Trilinos/BUILD_DIR" \
-D TPL_ENABLE_MPI:BOOL=ON \
-D CMAKE_C_FLAGS:STRING="-m64 -g" \
-D CMAKE_CXX_FLAGS:STRING="-m64 -g" \
-D CMAKE_Fortran_FLAGS:STRING="-m64 -g" \
-D Trilinos_ENABLE_ALL_PACKAGES:BOOL=OFF \
-D Trilinos_ENABLE_Zoltan:BOOL=ON \
-D Zoltan_ENABLE_EXAMPLES:BOOL=ON \
-D Zoltan_ENABLE_TESTS:BOOL=ON \
-D Zoltan_ENABLE_ParMETIS:BOOL=ON \
-D ParMETIS_INCLUDE_DIRS:FILEPATH="/home/username/code/ParMETIS3_1" \
-D ParMETIS_LIBRARY_DIRS:FILEPATH="/home/username/code/ParMETIS3_1" \
-D Zoltan_ENABLE_Scotch:BOOL=ON \
-D Scotch_INCLUDE_DIRS:FILEPATH="/home/username/code/scotch_5.1/include" \
-D Scotch_LIBRARY_DIRS:FILEPATH="/home/username/code/scotch_5.1/lib" \
..
make
make install
```

More examples are in the directory zoltan/SampleCmakeScripts. More details of CMake use in Trilinos are in Trilinos/cmake/TrilinosCMakeQuickstart.txt.

Testing the Zoltan Library

The *examples* directory contains simple C and C++ examples which use the Zoltan library. The makefile in this directory has three targets: These examples are built automatically when the [Autotools build environment](#) or [CMake build environment](#) is used.

The "right" answer for these tests depends on the number of processes with which you run the tests. In general, if they

compile successfully, run quickly (in seconds), and produce reasonable looking output, then Zoltan is built successfully.

Reporting Bugs in Zoltan

Zoltan uses [Bugzilla](#) to collect bug reports. Please read the [instructions for reporting bugs](#) through the Zoltan Bugzilla database.

Incorporating Zoltan into Applications

Incorporating Zoltan into applications requires three basic steps:

- Writing [query functions](#) that return information about the application to Zoltan.
- [Initializing](#) Zoltan, [creating a Zoltan object](#), and [setting appropriate parameters](#).
- Calling Zoltan tools to perform [partitioning](#), [ordering](#), [migration](#), [coloring](#), etc.

The set of [query functions](#) needed by an application depends on the particular tools (e.g., [partitioning](#), [ordering](#)) used and on the [algorithms](#) selected within the tools. Not all query functions are needed by every application. See documentation on tools and algorithms to determine which query functions are needed.

Building Applications that use Zoltan

The C library interface is described in the include file *include/zoltan.h*; this file should be included in all C application source files that call Zoltan library routines.

The [C++ interface](#) to Zoltan is implemented in header files which define classes that wrap the Zoltan C library. The file *include/zoltan_cpp.h* defines the **Zoltan** class which encapsulates a load balancing data structure and the Zoltan load balancing functions which operate upon it. Include this header file instead in your C++ application. Note that C++ applications should call the C function [Zoltan_Initialize](#) before creating a **Zoltan** object.

[Fortran applications](#) must USE [module zoltan](#) and specify the Zoltan installation's *include* directory as a directory to be searched for module information files.

The C, C++ or Fortran application should then be linked with the Zoltan library (built with Fortran support in the Fortran case) by including

-lzoltan

in the linking command for the application. Communication within Zoltan is performed through MPI, so appropriate MPI libraries must be linked with the application. Third-party libraries, such as [ParMETIS](#), [PT-Scotch](#) and [PaToH](#), must be also be linked with the application if they were included in compilation of the Zoltan library.

The installed files *include/Makefile.export.zoltan** contain macros that can specify Zoltan paths and libraries in an application's Makefiles. Using these files, applications can be assured they are using the same build options that were used when Zoltan was built.

Data Types for Object IDs

Application query functions and application callable library functions use global and local identifiers (IDs) for objects. *All objects to be used in load balancing must have unique global IDs.* Zoltan stores an ID as an array of unsigned

integers. The number of entries in these arrays can be set using the [NUM_GID_ENTRIES](#) and [NUM_LID_ENTRIES](#) parameters; by default, one unsigned integer represents an ID. Applications may use whatever format is most convenient to store their IDs; the IDs can then be converted to and from Zoltan's ID format in the [application-registered query functions](#).

The following type definitions are defined in *include/zoltan_types.h*; they can be used by an application for memory allocation, MPI communication, and as arguments to [load-balancing interface functions](#) and [application-registered query functions](#).

```
typedef unsigned int ZOLTAN_ID_TYPE;
typedef ZOLTAN_ID_TYPE *ZOLTAN_ID_PTR;
#define ZOLTAN_ID_MPI_TYPE MPI_UNSIGNED
```

In the Fortran interface, IDs are passed as arrays of integers since unsigned integers are not supported in Fortran. See the description of the [Fortran interface](#) for more details.

The local IDs passed to Zoltan are not used by the library; they are provided for the convenience of the application and can contain any information desired by the application. For instance, local array indices for objects may be passed as local IDs, enabling direct access to object data in the query function routines. See the [application-registered query functions](#) for more details. The source code distribution contains an example application *zdrive* in which global IDs are integers and local IDs are local array indices. One may choose not to use local ids at all, in which case [NUM_LID_ENTRIES](#) may be set to zero.

Some Zoltan routines (e.g., [Zoltan_LB_Partition](#) and [Zoltan_Invert_Lists](#)) allocate arrays of type **ZOLTAN_ID_PTR** and return them to the application. Others (e.g., [Zoltan_Order](#) and [Zoltan_DD_Find](#)) require the application to allocate memory for IDs. Memory for IDs can be allocated as follows:

```
ZOLTAN_ID_PTR gids;
int num_gids, int num_gid_entries;
gids = (ZOLTAN_ID_PTR) ZOLTAN\_MALLOC(num_gids * num_gid_entries *
sizeof(ZOLTAN_ID_TYPE));
```

The system call *malloc* may be used instead of [ZOLTAN_MALLOC](#).

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C++ Interface

The C++ interface to the Zoltan library is contained in the header files listed below. Each header file defines one class. Each class represents a Zoltan data structure and the functions that operate on that data structure. The class methods in the header files call functions in the Zoltan C library. So to use the C++ interface from your application, include the appropriate header file and link with the Zoltan C library.

header file	class
<i>include/zoltan_cpp.h</i>	Zoltan , representing a load balancing instance
<i>Utilities/Communication/zoltan_comm_cpp.h</i>	Zoltan_Comm , representing an unstructured communication instance
<i>Utilities/DDirectory/zoltan_dd_cpp.h</i>	Zoltan_DD , representing a distributed directory instance
<i>Utilities/Timer/zoltan_timer_cpp.h</i>	Zoltan_Timer , representing a timer instance

More detailed information about the interface may be found in the [Zoltan Developer's Guide](#).

Simple examples of the use of the interface may be found in the *examples/CPP* directory. A more complete example is the test driver [zCPPdrive](#). The source code for this test driver is in the *driver* directory.

A note on declaring application registered query functions from a C++ application may be found in the section titled [Application-Registered Query Functions](#).

Two peculiarities of the wrapping of Zoltan with C++ classes are mentioned here:

1. You must call the C language function [Zoltan_Initialize](#) before using the C++ interface to the Zoltan library. This function should only be called once. Due to design choices, the C++ interface maintains no global state that is independent of any instantiated objects, so it does not know if the function has been called or not. Therefore, the C++ wrappers do not call [Zoltan_Initialize](#) for you.
2. It is preferable to allocate **Zoltan** objects dynamically so you can explicitly delete them before your application exits. (**Zoltan** objects allocated instead on the stack will be deleted automatically at the completion of the scope in which they were created.) The reason is that the **Zoltan** destructor calls `Zoltan_Destroy()`, which makes an MPI call to free the communicator in use by the **Zoltan** object. However the MPI destructor may have been called before the **Zoltan** destructor. In this case you would receive an error while your application is exiting.

This second point is illustrated in the good and bad example below.

```
int main(int argc, char *argv[])
{
  MPI::Init(argc, argv);
  int rank = MPI::COMM_WORLD.Get_rank();
  int size = MPI::COMM_WORLD.Get_size();

  //Initialize the Zoltan library with a C language call
  float version;
  Zoltan_Initialize(argc, argv, &version);

  //Dynamically create Zoltan object.
  Zoltan *zz = new Zoltan(MPI::COMM_WORLD);
  zz->Set_Param("LB_METHOD", "RCB");

  //Several lines of code would follow, working with zz
```

```
//Explicitly delete the Zoltan object
delete zz;
MPI::Finalize();
}
```

Good example, Zoltan object is dynamically allocated and explicitly deleted before exit.

```
int main(int argc, char *argv[])
{
Zoltan zz;

MPI::Init(argc, argv);
int rank = MPI::COMM_WORLD.Get_rank();
int size = MPI::COMM_WORLD.Get_size();

//Initialize the Zoltan library with a C language call
float version;
Zoltan_Initialize(argc, argv, &version);

zz.Set_Param("LB_METHOD", "RCB");

//Several lines of code would follow, working with zz

MPI::Finalize();
}
```

Bad example, the MPI destructor may execute before the Zoltan destructor at process exit.

FORTRAN Interface

The Fortran interface for Zoltan is a Fortran 90 interface designed similar to the Fortran 90 Bindings for OpenGL [[Mitchell](#)]. There is no FORTRAN 77 interface; however, FORTRAN 77 applications can use Zoltan by adding only a few Fortran 90 statements, which are fully explained in the section on [FORTRAN 77](#), provided that vendor-specific extensions are not heavily used in the application. This section describes how to build the Fortran interface into the Zoltan library, how to call Zoltan from Fortran applications, and how to compile Fortran applications that use Zoltan. Note that the capitalization used in this section is for clarity and need not be adhered to in the application code, since Fortran is case insensitive.

- [Compiling Zoltan](#)
- [Compiling Applications](#)
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- [FORTRAN 77](#)
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FORTRAN: Compiling Zoltan

Including the Fortran interface in the Zoltan library requires an additional option on the [autotools configure](#) or [CMake](#) command:

- Autotools option: `--enable-f90interface`
- CMake option: `-D Zoltan_ENABLE_F90INTERFACE:BOOL=ON`

Before compiling the library, make sure that the application's [zoltan_user_data.f90](#) has been placed in the `zoltan/src/fort/` directory, if needed.

FORTRAN: Compiling Applications

To compile a Fortran application using the Zoltan library, the module information files must be made available to most compilers during the compilation phase. Module information files are files generated by the compiler to provide module information to program units that **USE** the module. They usually have suffixes like `.mod` or `.M`. The module information files for the modules in the Zoltan library are located in the *include* directory generated during [Zoltan installation](#). Most Fortran 90 compilers have a compile line flag to specify directories to be searched for module information files, typically `-I`; check the documentation for your compiler. If your compiler does not have such a flag, you will have to copy the module information files to the directory of the application (or use symbolic links).

The Fortran interface is built into the same library file as the rest of Zoltan, which is found during the compiler link phase with `-lzoltan`. Thus an example compilation line would be

```
f90 -I<path to the installation include directory> application.f90 -lzoltan
```

FORTRAN API

The Fortran interface for each [Zoltan Interface Function](#) and [Application-Registered Query Function](#) is given along with the C interface. This section contains some general information about the design and use of the Fortran interface.

- [Names](#)
- [Zoltan module](#)
- [Numeric types](#)
- [Structures](#)
- [Global and local IDs](#)
- [Query function data](#)

Names

All procedure, variable, defined constant and structure names are identical to those in the C interface, except that in Fortran they are case insensitive (either upper or lower case letters can be used).

Zoltan module

MODULE *zoltan* provides access to all entities in Zoltan that are of use to the application, including kind type parameters, named constants, procedures, and derived types. Any program unit (e.g., main program, module, external subroutine) that needs access to an entity from Zoltan must contain the statement

```
USE zoltan
```

near the beginning.

Numeric types

The correspondence between Fortran and C numeric types is achieved through the use of kind type parameters. In most cases, the default kind for a Fortran type will match the corresponding C type, but this is not guaranteed. To insure portability of the application code, it is highly recommended that the following kind type parameters be used in the declaration of all variables and constants that will be passed to a Zoltan procedure:

C	Fortran
int	INTEGER(KIND=Zoltan_INT)
float	REAL(KIND=Zoltan_FLOAT)
double	REAL(KIND=Zoltan_DOUBLE)

Note that "KIND=" is optional in declaration statements. The kind number for constants can be attached to the constant, e.g., 1.0_Zoltan_DOUBLE.

Structures

For any struct in the C interface to Zoltan, e.g. [Zoltan_Struct](#), there is a corresponding derived type in the Fortran interface. Variables of this type are declared as demonstrated below:

```
TYPE(Zoltan_Struct) :: zz
```

In the Fortran interface, the internal components of the derived type are PRIVATE and not accessible to the application. However, the application simply passes these variables around, and never needs to access the internal components.

Global and local IDs

While the C implementation uses arrays of unsigned integers to represent [global and local IDs](#), the Fortran interface uses arrays of integers, as unsigned integers are not available in Fortran. Thus, each ID is represented as an array (possibly of size 1) of integers. Applications that use other data types for their IDs can convert between their data types and Zoltan's in the [application-registered query functions](#).

Query function data

[Zoltan_Set_Fn](#) allows the application to pass a pointer to data that will subsequently be passed to the query function being registered. From Fortran this is an optional argument, or can be one of several types. In the simplest cases, an intrinsic array containing the data will be sufficient. For these cases, data can be an assumed size array of type INTEGER(Zoltan_INT), REAL(Zoltan_FLOAT) or REAL(Zoltan_DOUBLE). When the argument is omitted in the call to the registration function, a data argument will still be passed to the query function. This should be declared as an assumed size array of type INTEGER(Zoltan_INT) and never used.

For more complicated situations, the application may need to pass data in a user-defined type. The strong type checking of Fortran does not allow passing an arbitrary type without modifying the Fortran interface for each desired type. So the Fortran interface provides a type to be used for this purpose, **Zoltan_User_Data_1**. Since different types of data may need to be passed to different query functions, four such types are provided, using the numerals 1, 2, 3 and 4 as the last character in the name of the type. These types are defined by the application in *zoltan_user_data.f90*. If not needed, they must be defined, but can be almost empty as in *fort/zoltan_user_data.f90*.

The application may use these types in any appropriate way. If desired, it can define these types to contain the application's data and use the type throughout the application. But it is anticipated that in most cases, the desired type already exists in the application, and the **Zoltan_User_Data_x** types will be used as "wrapper types," containing one or more pointers to the existing types. For example,

```

TYPE mesh
    ! an existing data type with whatever defines a mesh
END TYPE mesh

TYPE Zoltan_User_Data_2

    TYPE(mesh), POINTER :: ptr
END TYPE Zoltan_User_Data_2
    
```

The application would then set the pointer to the data before calling Zoltan_Set_Fn:

```

TYPE(mesh) :: meshdata
TYPE(Zoltan_User_Data_2) :: query_data
TYPE(Zoltan_Struct) :: zz
INTEGER(Zoltan_INT), EXTERNAL :: num_obj_func ! not required for module procedures

query_data%ptr => meshdata
ierr = Zoltan_Set_Fn(zz,ZOLTAN_NUM_OBJ_FN_TYPE,num_obj_func,query_data)
    
```

Note that the existing data type must be available when **Zoltan_User_Data_x** is defined. Therefore it must be defined either in *zoltan_user_data.f90* or in a module that is compiled before *zoltan_user_data.f90* and **USED** by **MODULE** *zoltan_user_data*. For an example that uses a wrapper type, see *fdriver/zoltan_user_data.f90*.

FORTRAN 77

There is no FORTRAN 77 interface for Zoltan; however, an existing FORTRAN 77 application can be compiled by a

Fortran 90 compiler provided it does not use vendor specific extensions (unless the same extensions are supported by the Fortran 90 compiler), and the application can use Zoltan's Fortran 90 interface with a minimal amount of Fortran 90 additions. This section provides details of the Fortran 90 code that must be added.

When building the Zoltan library, use the file *fort/zoltan_user_data.f90* for *zoltan_user_data.f90*. This assumes that DATA in a call to [ZOLTAN_SET_FN](#) is either omitted (you can omit arguments that are labeled OPTIONAL in the Fortran API) or an array of type INTEGER, REAL or DOUBLE PRECISION (REAL*4 and REAL*8 might be acceptable). If a more complicated set of data is required (for example, two arrays), then it should be made available to the query functions through COMMON blocks.

To get access to the interface, each program unit (main program, subroutine or function) that calls a Zoltan routine must begin with the statement

```
USE ZOLTAN
```

and this should be the first statement after the program, subroutine or function statement (before the declarations).

The pointer to the Zoltan structure returned by [ZOLTAN_CREATE](#) should be declared as

```
TYPE(ZOLTAN_STRUCT), POINTER :: ZZ
```

(you can use a name other than ZZ if you wish).

To create the structure, use a pointer assignment statement with the call to [ZOLTAN_CREATE](#):

```
ZZ => ZOLTAN\_CREATE(COMMUNICATOR)
```

Note that the assignment operator is "=>". If ZZ is used in more than one procedure, then put it in a COMMON block. It cannot be passed as an argument unless the procedure interfaces are made "explicit." (Let's not go there.)

The eight import and export arrays passed to [ZOLTAN_LB_PARTITION](#) (and other procedures) must be pointers. They should be declared as, for example,

```
INTEGER, POINTER :: IMPORT_GLOBAL_IDS(:)
```

Note that the double colon after POINTER is required, and the dimension must be declared as "(:)" with a colon. Like ZZ, if they are used in more than one procedure, pass them through a COMMON block, not as an argument.

Except in the unlikely event that the default kinds of intrinsic types do not match the C intrinsic types, you do not have to use the kind type parameters **Zoltan_INT**, etc. It is also not necessary to include the INTENT attribute in the declarations of the query functions, so they can be simplified to, for example,

```
SUBROUTINE GET_OBJ_LIST(DATA, GLOBAL_IDS, LOCAL_IDS, WGT_DIM, OBJ_WGTS, IERR)
  INTEGER DATA(*),GLOBAL_IDS(*),LOCAL_IDS(*),WGT_DIM,IERR
  REAL OBJ_WGTS(*)
```

to be more consistent with a FORTRAN 77 style.

FORTRAN: System-Specific Remarks

System-specific details of the FORTRAN interface are included below.

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[MPICH](#)
[Pacific Sierra](#)
[NASoftware](#)

MPICH

As of version 1.1.2, the MPICH implementation of MPI is not completely "Fortran 90 friendly." Only one problem was encountered during our tests: the reliance on command line arguments. MPICH uses command line arguments during the start-up process, even if the application does not. Command line arguments are not standard in Fortran, so although most compilers offer it as an extension, each compiler has its own method of handling them. The problem arises when one Fortran compiler is specified during the build of MPICH and another Fortran compiler is used for the application. This should not be a problem on systems where there is only one Fortran compiler, or where multiple Fortran compilers are compatible (for example, FORTRAN 77 and Fortran 90 compilers from the same vendor). If your program can get past the call to MPI_Init, then you do not have this problem.

To solve this problem, build MPICH in such a way that it does not include the routines for *iargc* and *getarg* (I have been able to do this by using the -f95nag flag when configuring MPICH), and then provide your own versions of them when you link the application. Some versions of these routines are provided in *fdriver/farg_**.

Pacific Sierra

Pacific Sierra Research (PSR) Vastf90 is not currently supported due to bugs in the compiler with no known workarounds. It is not known when or if this compiler will be supported.

NASoftware

N.A.Software FortranPlus is not currently supported due to problems with the query functions. We anticipate that this problem can be overcome, and support will be added soon.

[\[Table of Contents\]](#) | [Next: Zoltan Interface Functions](#) | [Previous: C++ Interface](#) | [Privacy and Security](#)]

Zoltan Interface Functions

An application calls a series of dynamic load-balancing library functions to initialize the load balancer, perform load balancing and migrate data. This section describes the syntax of each type of interface function:

- [General Zoltan Interface Functions](#)
- [Load-Balancing Interface Functions](#)
- [Functions for Augmenting a Decomposition](#)
- [Migration Interface Functions](#)
- [Graph Coloring Functions](#)
- [Graph Ordering Functions](#)

Examples of the calling sequences for initialization, load-balancing, and data migration are included in the [Initialization](#), [Load-Balancing](#), and [Migration](#) sections, respectively, of the [Examples of Library Usage](#).

Error Codes

All interface functions, with the exception of [Zoltan_Create](#), return an error code to the application. The possible return codes are defined in `include/zoltan_types.h` and Fortran [module zoltan](#), and are listed in the [table](#) below.

Note: Robust error handling in parallel has not yet been achieved in Zoltan. When a processor returns from Zoltan due to an error condition, other processors do not necessarily return the same condition. In fact, other processors may not know that the original processor has returned from Zoltan, and may wait indefinitely in a communication routine (e.g., waiting for a message from the original processor that is not sent due to the error condition). The parallel error-handling capabilities of Zoltan will be improved in future releases.

<code>ZOLTAN_OK</code>	Function returned without warnings or errors.
<code>ZOLTAN_WARN</code>	Function returned with warnings. The application will probably be able to continue to run.
<code>ZOLTAN_FATAL</code>	A fatal error occured within the Zoltan library.
<code>ZOLTAN_MEMERR</code>	An error occurred while allocating memory. When this error occurs, the library frees any allocated memory and returns control to the application. If the application then wants to try to use another, less memory-intensive algorithm, it can do so.

Return codes defined in `include/zoltan_types.h`.

Naming conventions

The C, Fortran and C++ interfaces follow consistent naming conventions, as illustrated in the following table.

	C and Fortran	C++
Partitioning and migration functions example: perform partitioning example: assign a point to a part	<code>Zoltan_LB_function()</code> <code>Zoltan_LB_Partition()</code> <code>Zoltan_LB_Point_Assign()</code>	<code>Zoltan::function()</code> <code>Zoltan::LB_Partition()</code> <code>Zoltan::LB_Point_Assign()</code>

Unstructured communication example: perform communication	Zoltan_Comm <i>_function</i> Zoltan_Comm_Do()	Zoltan_Comm:: <i>function</i> Zoltan_Comm::Do()
Distributed data example: find objects in a remote process	Zoltan_DD <i>_function</i> Zoltan_DD_Find()	Zoltan_DD:: <i>function</i> Zoltan_DD::Find()
Timers example: print timing results	Zoltan_Timer <i>_function</i> Zoltan_Timer_Print()	Zoltan_Timer:: <i>function</i> Zoltan_Timer::Print()

In particular, the C++ **Zoltan** class represents a load balancing instance and the methods that operate on it. The method name is identical to the part of the C and Fortran function name that indicates the function performed. A C++ **Zoltan_Comm** object represents an instance of unstructured communication, a C++ **Zoltan_DD** object represents a distributed directory, and a C++ **Zoltan_Timer** object is a timer. Their method names are derived similarly.

General Interface Functions

Functions used to initialize and manipulate Zoltan's data structures are described below:

- [Zoltan_Initialize](#)
- [Zoltan_Create](#)
- [Zoltan_Copy](#)
- [Zoltan_Copy_To](#)
- [Zoltan_Set_Param](#)
- [Zoltan_Set_Param_Vec](#)
- [Zoltan_Set_Fn](#)
- [Zoltan_Set <zoltan_fn_type> Fn](#)
- [Zoltan_Destroy](#)

C and C++:	<pre>int Zoltan_Initialize (int <i>argc</i>, char **<i>argv</i>, float *<i>ver</i>);</pre>
FORTRAN:	<pre>FUNCTION Zoltan_Initialize(<i>argc</i>, <i>argv</i>, <i>ver</i>) INTEGER(Zoltan_INT) :: Zoltan_Initialize INTEGER(Zoltan_INT), INTENT(IN), OPTIONAL :: <i>argc</i> CHARACTER(LEN=*), DIMENSION(*), INTENT(IN), OPTIONAL :: <i>argv</i> REAL(Zoltan_FLOAT), INTENT(OUT) :: <i>ver</i></pre>

The **Zoltan_Initialize** function initializes MPI for Zoltan. If the application uses MPI, this function should be called after calling **MPI_Init**. If the application does not use MPI, this function calls **MPI_Init** for use by Zoltan. This function is called with the *argc* and *argv* command-line arguments from the main program, which are used if **Zoltan_Initialize** calls **MPI_Init**. From C, if **MPI_Init** has already been called, the *argc* and *argv* arguments may have any value because their values will be ignored. From Fortran, if one of *argc* or *argv* is omitted, they must both be omitted. If they are omitted, *ver* does NOT have to be passed as a keyword argument.

Zoltan_Initialize returns the Zoltan version number so that users can verify which version of the library their application is linked to.

C++ applications should call the C **Zoltan_Initialize** function before using the C++ interface to the Zoltan library.

Arguments:

- | | |
|-------------|-------------------------------------------------------------------------------|
| <i>argc</i> | The number of command-line arguments to the application. |
| <i>argv</i> | An array of strings containing the command-line arguments to the application. |
| <i>ver</i> | Upon return, the version number of the library. |

Returned Value:

- | | |
|-----|------------------------------|
| int | Error code . |
|-----|------------------------------|

C:	<pre>struct Zoltan_Struct *Zoltan_Create (MPI_Comm <i>communicator</i>);</pre>
FORTRAN:	<pre>FUNCTION Zoltan_Create(<i>communicator</i>)</pre>

```
TYPE(Zoltan_Struct), pointer :: Zoltan_Create
INTEGER, INTENT(IN) :: communicator

C++:      Zoltan (
           const MPI_Comm &communicator = MPI_COMM_WORLD);
```

The **Zoltan_Create** function allocates memory for storage of information to be used by Zoltan and sets the default values for the information. The pointer returned by this function is passed to many subsequent functions. An application may allocate more than one **Zoltan_Struct** data structure; for example, an application may use several **Zoltan_Struct** structures if, say, it uses different decompositions with different load-balancing techniques.

In the C++ interface to Zoltan, the **Zoltan** class represents a Zoltan load balancing data structure and the functions that operate on it. It is the constructor which allocates an instance of a **Zoltan** object. It has no return value.

Arguments:

communicator The MPI communicator to be used for this Zoltan structure. Only those processors included in the communicator participate in Zoltan functions. If all processors are to participate, *communicator* should be **MPI_COMM_WORLD** .

Returned Value:

struct **Zoltan_Struct** Pointer to memory for storage of Zoltan information. If an error occurs, NULL will be returned in C, or the result will be a nullified pointer in Fortran. Any error that occurs in this function is assumed to be fatal.

```
C:      struct Zoltan_Struct *Zoltan_Copy (
           Zoltan_Struct *from);

FORTRAN:  FUNCTION Zoltan_Copy(from)
           TYPE(Zoltan_Struct), pointer :: Zoltan_Copy
           TYPE(Zoltan_Struct), INTENT(IN) :: from

C++:      Zoltan (
           const Zoltan &zz);
```

The **Zoltan_Copy** function creates a new **Zoltan_Struct** and copies the state of the existing **Zoltan_Struct**, which it has been passed, to the new structure. It returns the new **Zoltan_Struct**.

There is no direct interface to **Zoltan_Copy** from C++. Rather, the **Zoltan** copy constructor invokes the C library **Zoltan_Copy** program.

Arguments:

from A pointer to the **Zoltan_Struct** that is to be copied.

Returned Value:

struct **Zoltan_Struct** Pointer to a new **Zoltan_Struct**, which is now a copy of *from*.

```
C:      int Zoltan_Copy_To (
           Zoltan_Struct *to,
           Zoltan_Struct *from);

FORTRAN:  FUNCTION Zoltan_Copy_To(to, from)
           INTEGER(Zoltan_INT) :: Zoltan_Copy_To
```

```
TYPE(Zoltan_Struct), INTENT(IN) :: to
TYPE(Zoltan_Struct), INTENT(IN) :: from

C++:      Zoltan & operator= (
           const Zoltan &zz);
```

The **Zoltan_Copy_To** function copies one **Zoltan_Struct** to another, after first freeing any memory used by the target **Zoltan_Struct** and re-initializing it.

The C++ interface to the **Zoltan_Copy_To** function is through the **Zoltan** copy operator, which invokes the C library **Zoltan_Copy_To** program.

Arguments:

<i>to</i>	A pointer to an existing Zoltan_Struct , the target of the copy.
<i>from</i>	A pointer to an existing Zoltan_Struct , the source of the copy.

Returned Value:

int	0 on success and 1 on failure.
-----	----------------------------------------------

```
C:      int Zoltan_Set_Param (
        struct Zoltan_Struct *zz,
        char *param_name,
        char *new_val);

FORTRAN: FUNCTION Zoltan_Set_Param(zz, param_name, new_val)
          INTEGER(Zoltan_INT) :: Zoltan_Set_Param
          TYPE(Zoltan_Struct), INTENT(IN) :: zz
          CHARACTER(LEN=*), INTENT(IN) :: param_name, new_value

C++:    int Zoltan::Set_Param (
        const std::string &param_name,
        const std::string &new_value);
```

Zoltan_Set_Param is used to alter the value of one of the parameters used by Zoltan. All Zoltan parameters have reasonable default values, but this routine allows a user to provide alternative values if desired.

Arguments:

<i>zz</i>	Pointer to the Zoltan structure created by Zoltan_Create .
<i>param_name</i>	A string containing the name of the parameter to be altered. Note that the string is case-insensitive. Also, different Zoltan structures can have different parameter values.
<i>new_val</i>	A string containing the new value for the parameter. Example strings include "3.154", "True", "7" or anything appropriate for the parameter being set. As above, the string is case-insensitive.

Returned Value:

int	Error code .
-----	------------------------------

```
C:      int Zoltan_Set_Param_Vec (
        struct Zoltan_Struct *zz,
        char *param_name,
        char *new_val,
        int index);
```



```
FORTRAN:      FUNCTION Zoltan_Set_Param_Vec(zz, param_name, new_val, index)
               INTEGER(Zoltan_INT) :: Zoltan_Set_Param_Vec
               TYPE(Zoltan_Struct), INTENT(IN) :: zz
               CHARACTER(LEN=*), INTENT(IN) :: param_name, new_value
               INTEGER(Zoltan_INT), INTENT(IN) :: index

C++:          int Zoltan::Set_Param_Vec (
               const std::string &param_name,
               const std::string &new_val,
               const int &index);
```

Zoltan_Set_Param_Vec is used to alter the value of a vector parameter in Zoltan. A vector parameter is a parameter that has one name but contains multiple values. These values are referenced by their indices, usually starting at 0. Each entry (component) may have a different value. This routine sets a single entry (component) of a vector parameter. If you want all entries (components) of a vector parameter to have the same value, set the parameter using [Zoltan_Set_Param](#) as if it were a scalar parameter. If one only sets the values of a subset of the indices for a vector parameter, the remaining entries will have the default value for that particular parameter.

Arguments:

<i>zz</i>	Pointer to the Zoltan structure created by Zoltan_Create .
<i>param_name</i>	A string containing the name of the parameter to be altered. Note that the string is case-insensitive. Also, different Zoltan structures can have different parameter values.
<i>new_val</i>	A string containing the new value for the parameter. Example strings include "3.154", "True", "7" or anything appropriate for the parameter being set. As above, the string is case-insensitive.
<i>index</i>	The index of the entry of the vector parameter to be set. The default in Zoltan is that the first entry in a vector has index 0 (C-style indexing).

Returned Value:

int	Error code .
-----	------------------------------

```
C:            int Zoltan_Set_Fn (
               struct Zoltan_Struct *zz,
               ZOLTAN_FN_TYPE fn_type,
               void (*fn_ptr)(),
               void *data);

FORTRAN:      FUNCTION Zoltan_Set_Fn(zz, fn_type, fn_ptr, data)
               INTEGER(Zoltan_INT) :: Zoltan_Set_Fn
               TYPE(Zoltan_Struct), INTENT(IN) :: zz
               TYPE(ZOLTAN_FN_TYPE), INTENT(IN) :: fn_type
               EXTERNAL :: fn_ptr
               <type-data>, OPTIONAL :: data

               <type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),
               DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x)
               where x is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.

C++:          int Zoltan::Set_Fn (
               const ZOLTAN_FN_TYPE &fn_type,
               void (*fn_ptr)(),
               void *data = 0);
```

Zoltan_Set_Fn registers an application-supplied query function in the Zoltan structure. All types of query functions

can be registered through calls to **Zoltan_Set_Fn**. To register functions while maintaining strict type-checking of the *fn_ptr* argument, use [Zoltan_Set_<zoltan_fn_type>_Fn](#).

Arguments:

<i>zz</i>	Pointer to the Zoltan structure created by Zoltan_Create .
<i>fn_type</i>	The type of function being registered; see Application-Registered Query Functions for possible function types.
<i>fn_ptr</i>	A pointer to the application-supplied query function being registered.
<i>data</i>	A pointer to user defined data that will be passed, as an argument, to the function pointed to by <i>fn_ptr</i> . In C it may be NULL. In Fortran it may be omitted.

Returned Value:

int	Error code .
-----	------------------------------

```
C:      int Zoltan_Set_<zoltan_fn_type>_Fn (
        struct Zoltan_Struct *zz,
        <zoltan_fn_type> (*fn_ptr)(),
        void *data);

FORTRAN: FUNCTION Zoltan_Set_<zoltan_fn_type>_Fn(zz, fn_ptr, data)
          INTEGER(Zoltan_INT) :: Zoltan_Set_<zoltan_fn_type>_Fn
          TYPE(Zoltan_Struct), INTENT(IN) :: zz
          EXTERNAL :: fn_ptr
          <type-data>, OPTIONAL :: data
```

An interface block for *fn_ptr* is included in the FUNCTION definition so that strict type-checking of the registered query function can be done.

<type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

```
C++:    int Zoltan::Set_<zoltan_fn_type>_Fn (
        <zoltan_fn_type> (*fn_ptr)(),
        void *data = 0);
```

The interface functions **Zoltan_Set_<zoltan_fn_type>_Fn**, where [<zoltan_fn_type>](#) is one of the query function types, register specific types of [application-supplied query functions](#) in the Zoltan structure. One interface function exists for each type of query function. For example, **Zoltan_Set_Num_Geom_Fn** registers a query function of type [ZOLTAN_NUM_GEOM_FN](#). Each query function has an associated **Zoltan_Set_<zoltan_fn_type>_Fn**. A complete list of these functions is included in *include/zoltan.h*.

Query functions can be registered using either [Zoltan_Set_Fn](#) or **Zoltan_Set_<zoltan_fn_type>_Fn**. **Zoltan_Set_<zoltan_fn_type>_Fn** provides strict type checking of the *fn_ptr* argument; the argument's type is specified for each **Zoltan_Set_<zoltan_fn_type>_Fn**. [Zoltan_Set_Fn](#) does not provide this strict type checking, as the pointer to the registered function is cast to a void pointer.

Arguments:

<i>zz</i>	Pointer to the Zoltan structure created by Zoltan_Create .
<i>fn_ptr</i>	A pointer to the application-supplied query function being registered. The type of the pointer matches <zoltan_fn_type> in the name Zoltan_Set_<zoltan_fn_type>_Fn .
<i>data</i>	A pointer to user defined data that will be passed, as an argument, to the function pointed to by <i>fn_ptr</i> . In C it may be NULL. In Fortran it may be omitted.

Returned Value:

int [Error code](#).

Example:

The interface function
int **Zoltan_Set_Geom_Fn**(struct **Zoltan_Struct** *zz, [ZOLTAN_GEOM_FN](#) (*fn_ptr)(),
 void *data);
registers an [ZOLTAN_GEOM_FN](#) query function.

C:	void Zoltan_Destroy (struct Zoltan_Struct **zz);
FORTTRAN:	SUBROUTINE Zoltan_Destroy (zz) TYPE(Zoltan_Struct), POINTER :: zz
C++:	~ Zoltan ();

Zoltan_Destroy frees the memory associated with a Zoltan structure and sets the structure to NULL in C or nullifies the structure in Fortran. Note that **Zoltan_Destroy** does not deallocate the import and export arrays returned from Zoltan (e.g., the arrays returned from [Zoltan_LB_Partition](#)); these arrays can be deallocated through a separate call to [Zoltan_LB_Free_Part](#).

There is no explicit **Destroy** method in the C++ interface. The **Zoltan** object is destroyed when the destructor executes.

As a side effect, **Zoltan_Destroy** (and the C++ **Zoltan** destructor) frees the MPI communicator that had been allocated for the structure. So it is important that the application does not call **MPI_Finalize** before it calls **Zoltan_Destroy** or before the destructor executes.

Arguments:

zz A pointer to the address of the Zoltan structure, created by [Zoltan_Create](#), to be destroyed.

Load-Balancing Functions

The following functions are the load-balancing interface functions in the Zoltan library; their descriptions are included below.

[Zoltan_LB_Partition](#)

[Zoltan_LB_Set_Part_Sizes](#)

[Zoltan_LB_Eval](#)

[Zoltan_LB_Free_Part](#)

For [backward compatibility](#) with previous versions of Zoltan, the following functions are also maintained. These functions are applicable only when the number of parts to be generated is equal to the number of processors on which the parts are computed. That is, these functions assume "parts" and "processors" are synonymous.

[Zoltan_LB_Balance](#)

[Zoltan_LB_Free_Data](#)

Descriptions of algorithm-specific interface functions are included with the documentation of their associated algorithms. Algorithm-specific functions include:

[Zoltan_RCB_Box](#)

```

C:          int Zoltan_LB_Partition (
              struct Zoltan_Struct *zz,
              int *changes,
              int *num_gid_entries,
              int *num_lid_entries,
              int *num_import,
              ZOLTAN\_ID\_PTR *import_global_ids,
              ZOLTAN\_ID\_PTR *import_local_ids,
              int **import_procs,
              int **import_to_part,
              int *num_export,
              ZOLTAN\_ID\_PTR *export_global_ids,
              ZOLTAN\_ID\_PTR *export_local_ids,
              int **export_procs,
              int **export_to_part);

FORTRAN:    FUNCTION Zoltan_LB_Partition(zz, changes, num_gid_entries, num_lid_entries, num_import,
import_global_ids, import_local_ids, import_procs, import_to_part, num_export,
export_global_ids, export_local_ids, export_procs, export_to_part)
INTEGER(Zoltan_INT) :: Zoltan_LB_Partition
TYPE(Zoltan_Struct), INTENT(IN) :: zz
LOGICAL, INTENT(OUT) :: changes
INTEGER(Zoltan_INT), INTENT(OUT) :: num_gid_entries, num_lid_entries
INTEGER(Zoltan_INT), INTENT(OUT) :: num_import, num_export
INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_global_ids, export_global_ids
INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_local_ids, export_local_ids
INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_procs, export_procs
INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_to_part, export_to_part

```

```
C++:      int Zoltan::LB_Partition (
            int &changes,
            int &num_gid_entries,
            int &num_lid_entries,
            int &num_import,
            ZOLTAN\_ID\_PTR &import_global_ids,
            ZOLTAN\_ID\_PTR &import_local_ids,
            int * &import_procs,
            int * &import_to_part,
            int &num_export,
            ZOLTAN\_ID\_PTR &export_global_ids,
            ZOLTAN\_ID\_PTR &export_local_ids,
            int * &export_procs,
            int * &export_to_part);
```

Zoltan_LB_Partition invokes the load-balancing routine specified by the [LB_METHOD](#) parameter. The number of parts it generates is specified by the [NUM_GLOBAL_PARTS](#) or [NUM_LOCAL_PARTS](#) parameters. Results of the partitioning are returned in lists of objects to be imported into and exported from parts on this processor. Objects are included in these lists if *either* their part assignment or their processor assignment is changed by the new decomposition. If an application requests multiple parts on a single processor, these lists may include objects whose part assignment is changing, but whose processor assignment is unchanged.

Returned arrays are allocated in Zoltan; applications should not allocate these arrays before calling **Zoltan_LB_Partition**. The arrays are later freed through calls to [Zoltan_LB_Free_Part](#).

Arguments:

<i>zz</i>	Pointer to the Zoltan structure, created by Zoltan_Create , to be used in this invocation of the load-balancing routine.
<i>changes</i>	Set to 1 or .TRUE. if the decomposition was changed by the load-balancing method; 0 or .FALSE. otherwise.
<i>num_gid_entries</i>	Upon return, the number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	Upon return, the number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES .
<i>num_import</i>	Upon return, the number of objects that are newly assigned to this processor or to parts on this processor (i.e., the number of objects being imported from different parts to parts on this processor). If the value returned is -1, no import information has been returned and all import arrays below are NULL. (The RETURN_LISTS parameter determines whether import lists are returned).
<i>import_global_ids</i>	Upon return, an array of <i>num_import</i> global IDs of objects to be imported to parts on this processor. (size = <i>num_import</i> * <i>num_gid_entries</i>)
<i>import_local_ids</i>	Upon return, an array of <i>num_import</i> local IDs of objects to be imported to parts on this processor. (size = <i>num_import</i> * <i>num_lid_entries</i>)
<i>import_procs</i>	Upon return, an array of size <i>num_import</i> listing the processor IDs of the processors that owned the imported objects in the previous decomposition (i.e., the source processors).
<i>import_to_part</i>	Upon return, an array of size <i>num_import</i> listing the parts to which the imported objects are being imported.
<i>num_export</i>	Upon return, this value of this count and the following lists depends on the value of the RETURN_LISTS parameter:

- It is the count of objects on this processor that are newly assigned to other processors or to other parts on this processor, if [RETURN_LISTS](#) is "EXPORT" or "EXPORT AND IMPORT".
- It is the count of all objects on this processor, if [RETURN_LISTS](#) is "PARTS" (or "PART ASSIGNMENTS").
- It is -1 if the value of [RETURN_LISTS](#) indicates that either no lists are to be returned, or only import lists are to be returned. If the value returned is -1, no export information has been returned and all export arrays below are NULL .

<i>export_global_ids</i>	Upon return, an array of <i>num_export</i> global IDs of objects to be exported from parts on this processor (if RETURN_LISTS is equal to "EXPORT" or "EXPORT AND IMPORT"), or an array of <i>num_export</i> global IDs for every object on this processor (if RETURN_LISTS is equal to "PARTS" or "PART ASSIGNMENTS"), . (size = <i>num_export</i> * <i>num_gid_entries</i>)
<i>export_local_ids</i>	Upon return, an array of <i>num_export</i> local IDs associated with the global IDs returned in <i>export_global_ids</i> (size = <i>num_export</i> * <i>num_lid_entries</i>)
<i>export_procs</i>	Upon return, an array of size <i>num_export</i> listing the processor ID of the processor to which each object is now assigned (i.e., the destination processor). If RETURN_LISTS is equal to "PARTS" or "PART ASSIGNMENTS", this list includes all objects, otherwise it only includes the objects which are moving to a new part and/or process.
<i>export_to_part</i>	Upon return, an array of size <i>num_export</i> listing the parts to which the objects are assigned under the new partition.

Returned Value:

int [Error code](#).

```
C:      int Zoltan_LB_Set_Part_Sizes (  
        struct Zoltan_Struct *zz,  
        int global_num,  
        int len,  
        int *part_ids,  
        int *wgt_idx,  
        float *part_sizes);  
  
FORTRAN: function Zoltan_LB_Set_Part_Sizes( zz,global_part,len,partids,wgtidx,partsizes)  
integer(Zoltan_INT) :: Zoltan_LB_Set_Part_Sizes  
type(Zoltan_Struct) INTENT(IN) zz  
integer(Zoltan_INT) INTENT(IN) global_part,len,partids(*),wgtidx(*)  
real(Zoltan_FLOAT) INTENT(IN) partsizes(*)  
  
C++:    int Zoltan::LB_Set_Part_Sizes (  
        const int &global_num,  
        const int &len,  
        int *part_ids,  
        int *wgt_idx,  
        float *part_sizes);
```

Zoltan_LB_Set_Part_Sizes is used to specify the desired part sizes in Zoltan. By default, Zoltan assumes that all parts should be of equal size. With **Zoltan_LB_Set_Part_Sizes**, one can specify the relative (not absolute) sizes of the parts. For example, if two parts are requested and the desired sizes are 1 and 2, that means that the first part will be assigned approximately one third of the total load. If the sizes were instead given as 1/3 and 2/3, respectively, the result would be exactly the same. Note that if there are multiple weights per object, one can (must) specify the part size

for each weight dimension independently.

Arguments:

<i>zz</i>	Pointer to the Zoltan structure created by Zoltan_Create .
<i>global_num</i>	Set to 1 if global part numbers are given, 0 otherwise (local part numbers).
<i>len</i>	Length of the next three input arrays.
<i>part_ids</i>	Array of part numbers, either global or local. (Part numbers are integers starting from 0.)
<i>vwgt_idx</i>	Array of weight indices (between 0 and OBJ_WEIGHT_DIM-1). This array should contain all zeros when there is only one weight per object.
<i>part_sizes</i>	Relative values for part sizes; <i>part_sizes[i]</i> is the desired relative size of the <i>vwgt_idx[i]</i> 'th weight of part <i>part_ids[i]</i> .

Returned Value:

int	Error code .
-----	------------------------------

```
C:      int Zoltan_LB_Eval (  
        struct Zoltan_Struct *zz,  
        int print_stats,  
        ZOLTAN_BALANCE_EVAL *obj_info,  
        ZOLTAN_GRAPH_EVAL *graph_info,  
        ZOLTAN_HG_EVAL *hg_info);  
  
FORTRAN:  FUNCTION Zoltan_LB_Eval(zz, print_stats)  
          INTEGER(Zoltan_INT) :: Zoltan_LB_Eval  
          TYPE(Zoltan_Struct), INTENT(IN) :: zz  
          LOGICAL, INTENT(IN) :: print_stats  
  
C++:      int Zoltan::LB_Eval (  
        const int &print_stats,  
        ZOLTAN_BALANCE_EVAL *obj_info,  
        ZOLTAN_GRAPH_EVAL *graph_info,  
        ZOLTAN_HG_EVAL *hg_info);
```

Zoltan_LB_Eval evaluates the quality of a decomposition. The quality metrics of interest differ depending on how you are using Zoltan. If you are partitioning points in space using one of Zoltan's geometric methods, you will want to know the weighted balance of objects across parts. However if you are partitioning a graph, you will want to know about edges that have vertices in more than one part. **Zoltan_LB_Eval** can write three different structures with these differing metrics.

Arguments:

<i>zz</i>	Pointer to the Zoltan structure.
<i>print_stats</i>	If <i>print_stats</i> >0 (.TRUE. in Fortran), print the quality metrics to <i>stdout</i> .
<i>obj_info</i>	If <i>obj_info</i> is non-NULL, write object balance values to the ZOLTAN_BALANCE_EVAL structure.
<i>graph_info</i>	If <i>graph_info</i> is non-NULL, write graph partition metrics to the ZOLTAN_GRAPH_EVAL structure.
<i>hg_info</i>	If <i>hg_info</i> is non-NULL, write hypergraph partition metrics to the ZOLTAN_HG_EVAL structure.

Returned Value:

int	Error code .
-----	------------------------------

The EVAL structures are defined in `zoltan/src/include/zoltan_eval.h`. Several of the fields in the EVAL structures are arrays of values. The arrays contain values for

1. the total for the local process
2. the total across all parts
3. the minimum across all parts
4. the maximum across all parts
5. the average across all parts

in that order. The corresponding macros that refer to these fields are:

```
#define EVAL_LOCAL_SUM 0
#define EVAL_GLOBAL_SUM 1
#define EVAL_GLOBAL_MIN 2
#define EVAL_GLOBAL_MAX 3
#define EVAL_GLOBAL_AVG 4
```

The ZOLTAN_BALANCE_EVAL structure contains the following fields, and would be of interest if you are doing geometric partitioning:

- **obj_imbalance**: imbalance in count of objects across parts, scaled by requested part sizes.
- **imbalance**: imbalance in weight of objects across parts, scaled by requested part sizes.
- **nobj**: an array containing the number of objects on the local *process*, the total number of objects, the minimum number of objects in a *part*, the maximum number in a partition, and the average number of objects across the parts.
- **obj_wgt**: an array containing the sum of the weight of objects on the local *process*, the total weight across all processes, the minimum weight in a *part*, the maximum weight in a partition, and the average weight of objects across the parts.
- **xtra_imbalance**: if the [OBJ_WEIGHT_DIM](#) exceeds one, the **obj_imbalance** value for the extra weights is in this array.
- **xtra_obj_wgt**: if the [OBJ_WEIGHT_DIM](#) exceeds one, the **obj_wgt** array for the extra weights is in this array.

The ZOLTAN_GRAPH_EVAL structure contains the following fields, and would be of interest if you are doing graph partitioning:

- **cutl**: an array containing what is known at the CUTL or the ConCut measure of the graph, the first element is not set (this is the local process total, which has no meaning), the next is the sum of the CUTL across parts, then the minimum CUTL across parts, then the maximum CUTL across parts, finally the average CUTL across parts.
- **cutn**: an array containing what is known at the CUTN or the NetCut measure of the graph, the first element is not set, the next is the sum of the CUTL across parts, then the minimum CUTL across parts, then the maximum CUTL across parts, finally the average CUTL across parts.
- **cuts**: an array counting the number of cut edges (the first element again is not set) across all parts, the minimum across parts of cut edges, the maximum, and then the average
- **cut_weight**: an array summing the weight of cut edges (the first element again is not set) across all parts, the minimum across parts, the maximum, and then the average
- **nborparts**: an array which counts the number of neighboring parts that each part has, the first element again is not set, the next is the sum across all parts, the minimum across parts, the maximum, and then the average
- **obj_imbalance**: imbalance in count of objects across parts, scaled by requested part sizes.
- **imbalance**: imbalance in weight of objects across parts, scaled by requested part sizes.
- **nobj**: an array containing the number of objects on the local *process*, the total number of objects, the minimum number of objects in a *part*, the maximum number in a partition, and the average number of objects across the parts.
- **obj_wgt**: an array containing the sum of the weight of objects on the local *process*, the total weight across all processes, the minimum weight in a *part*, the maximum weight in a partition, and the average weight of objects across the parts.
- **num_boundary**: an array which counts the number of objects in a part that have at least one remote neighbor,

the first element is not set, the next is the sum across parts, the next is the minimum count in any part, the next is the maximum, and then the average for all parts

- **xtra_imbalance**: if the [OBJ_WEIGHT_DIM](#) exceeds one, the **obj_imbalance** value for the extra weights is in this array.
- **xtra_obj_wgt**: if the [OBJ_WEIGHT_DIM](#) exceeds one, the **obj_wgt** array for the extra weights is in this array.
- **xtra_cut_wgt**: if the [EDGE_WEIGHT_DIM](#) exceeds one, the **cut_wgt** array for the each extra weight is in this array.

The ZOLTAN_HG_EVAL structure contains the following fields, and would be of interest if you are doing hypergraph partitioning:

- **obj_imbalance**: imbalance in count of objects across parts, scaled by requested part sizes.
- **imbalance**: imbalance in weight of objects across parts, scaled by requested part sizes.
- **cutl**: an array containing what is known at the CUTL or the ConCut measure of the hypergraph, the first element is not set (this is the local process total, which has no meaning), the next is the sum of the CUTL across parts, then the minimum CUTL across parts, then the maximum CUTL across parts, finally the average CUTL across parts.
- **cutn**: an array containing what is known at the CUTN or the NetCut measure of the hypergraph, the first element is not set, the next is the sum of the CUTL across parts, then the minimum CUTL across parts, then the maximum CUTL across parts, finally the average CUTL across parts.
- **nobj**: an array containing the number of objects on the local *process*, the total number of objects, the minimum number of objects in a *part*, the maximum number in a partition, and the average number of objects across the parts.
- **obj_wgt**: an array containing the sum of the weight of objects on the local *process*, the total weight across all processes, the minimum weight in a *part*, the maximum weight in a part, and the average weight of objects across the parts.
- **xtra_imbalance**: if the [OBJ_WEIGHT_DIM](#) exceeds one, the **obj_imbalance** value for the extra weights is in this array.
- **xtra_obj_wgt**: if the [OBJ_WEIGHT_DIM](#) exceeds one, the **obj_wgt** array for the extra weights is in this array.

**Query
functions:**

Required: [ZOLTAN_NUM_OBJ_FN](#) and [ZOLTAN_OBJ_LIST_FN](#)

[Graph-Based](#) functions are required for writing the [ZOLTAN_GRAPH_EVAL](#) structure.

Optional: [Hypergraph-Based](#) functions will be used for writing the [ZOLTAN_HG_EVAL](#) if they are available, otherwise Zoltan will create a hypergraph from the [graph-Based](#) functions.
If [Zoltan_LB_Set_Part_Sizes](#) has been called, the part sizes set by the [Zoltan_LB_Set_Part_Sizes](#) will be used in calculating imbalances.

Note that the sum of *ncuts* over all processors is actually twice the number of edges cut in the graph (because each edge is counted twice). The same principle holds for *cut_wgt*.

There are a few improvements in Zoltan_LB_Eval in Zoltan version 1.5 (or higher). First, the balance data are computed with respect to both processors and parts (if applicable). Second, the desired part sizes (as set by Zoltan_LB_Set_Part_Sizes) are taken into account when computing the imbalance.

Known bug: If a part is spread across several processors, the computed cut information (*ncuts* and *cut_wgt*) may be incorrect (too high).

C: int Zoltan_LB_Free_Part (

```
ZOLTAN\_ID\_PTR *global_ids,  
ZOLTAN\_ID\_PTR *local_ids,  
int **procs,  
int **to_part);
```

FORTTRAN: FUNCTION **Zoltan_LB_Free_Part**(*global_ids, local_ids, procs, to_part*)
 INTEGER(Zoltan_INT) :: Zoltan_LB_Free_Part
 INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: global_ids
 INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: local_ids
 INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: procs, to_part

C++: int **Zoltan::LB_Free_Part** (
 [ZOLTAN_ID_PTR](#) *global_ids,
 [ZOLTAN_ID_PTR](#) *local_ids,
 int **procs,
 int **to_part);

Zoltan_LB_Free_Part frees the memory allocated by Zoltan to return the results of [Zoltan_LB_Partition](#) or [Zoltan_Invert_Lists](#). Memory pointed to by the arguments is freed and the arguments are set to NULL in C and C++ or nullified in Fortran. NULL arguments may be passed to **Zoltan_LB_Free_Part**. Typically, **Zoltan_LB_Free_Part** is called twice: once for the import lists, and once for the export lists. Note that this function does not destroy the Zoltan data structure itself; it is deallocated through a call to [Zoltan_Destroy](#) in C and Fortran and by the object destructor in C++.

Arguments:

<i>global_ids</i>	An array containing the global IDs of objects.
<i>local_ids</i>	An array containing the local IDs of objects.
<i>procs</i>	An array containing processor IDs.
<i>to_part</i>	An array containing part numbers.

Returned Value:

int	Error code .
-----	------------------------------

C: int **Zoltan_LB_Balance** (
 struct **Zoltan_Struct** *zz,
 int *changes,
 int *num_gid_entries,
 int *num_lid_entries,
 int *num_import,
 [ZOLTAN_ID_PTR](#) *import_global_ids,
 [ZOLTAN_ID_PTR](#) *import_local_ids,
 int **import_procs,
 int *num_export,
 [ZOLTAN_ID_PTR](#) *export_global_ids,
 [ZOLTAN_ID_PTR](#) *export_local_ids,
 int **export_procs);

FORTTRAN: FUNCTION **Zoltan_LB_Balance**(*zz, changes, num_gid_entries, num_lid_entries, num_import, import_global_ids, import_local_ids, import_procs, num_export, export_global_ids, export_local_ids, export_procs*)
 INTEGER(Zoltan_INT) :: Zoltan_LB_Balance
 TYPE(Zoltan_Struct), INTENT(IN) :: zz
 LOGICAL, INTENT(OUT) :: changes
 INTEGER(Zoltan_INT), INTENT(OUT) :: num_gid_entries, num_lid_entries
 INTEGER(Zoltan_INT), INTENT(OUT) :: num_import, num_export

INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_global_ids, export_global_ids
INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_local_ids, export_local_ids
INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_procs, export_procs

Zoltan_LB_Balance is a wrapper around [Zoltan_LB_Partition](#) that excludes the part assignment results. **Zoltan_LB_Balance** assumes the number of parts is equal to the number of processors; thus, the part assignment is equivalent to the processor assignment. Results of the partitioning are returned in lists of objects to be imported and exported. These arrays are allocated in Zoltan; applications should not allocate these arrays before calling **Zoltan_LB_Balance**. The arrays are later freed through calls to [Zoltan_LB_Free_Data](#) or [Zoltan_LB_Free_Part](#).

Arguments:

All arguments are analogous to those in [Zoltan_LB_Partition](#). Part-assignment arguments *import_to_part* and *export_to_part* are not included, as processor and parts numbers are considered to be the same in **Zoltan_LB_Balance**.

Returned Value:

int [Error code](#).

```
C:      int Zoltan_LB_Free_Data (
          ZOLTAN\_ID\_PTR *import_global_ids,
          ZOLTAN\_ID\_PTR *import_local_ids,
          int **import_procs,
          ZOLTAN\_ID\_PTR *export_global_ids,
          ZOLTAN\_ID\_PTR *export_local_ids,
          int **export_procs);

FORTRAN: FUNCTION Zoltan_LB_Free_Data(import_global_ids, import_local_ids, import_procs,
                                     export_global_ids, export_local_ids, export_procs)
          INTEGER(Zoltan_INT) :: Zoltan_LB_Free_Data
          INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_global_ids, export_global_ids
          INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_local_ids, export_local_ids
          INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_procs, export_procs
```

Zoltan_LB_Free_Data frees the memory allocated by the Zoltan to return the results of [Zoltan_LB_Balance](#) or [Zoltan_Compute_Destinations](#). Memory pointed to by the arguments is freed and the arguments are set to NULL in C or nullified in Fortran. NULL arguments may be passed to **Zoltan_LB_Free_Data**. Note that this function does not destroy the Zoltan data structure itself; it is deallocated through a call to [Zoltan_Destroy](#).

Arguments:

- | | |
|--------------------------|-------------------------------------------------------------------------------------------------------------------------------------------------------|
| <i>import_global_ids</i> | The array containing the global IDs of objects imported to this processor. |
| <i>import_local_ids</i> | The array containing the local IDs of objects imported to this processor. |
| <i>import_procs</i> | The array containing the processor IDs of the processors that owned the imported objects in the previous decomposition (i.e., the source processors). |
| <i>export_global_ids</i> | The array containing the global IDs of objects exported from this processor. |
| <i>export_local_ids</i> | The array containing the local IDs of objects exported from this processor. |
| <i>export_procs</i> | The array containing the processor IDs of processors that own the exported objects in the new decomposition (i.e., the destination processors). |

Returned Value:

int [Error code](#).

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Functions for Augmenting a Decomposition

The following functions support the addition of new items to an existing decomposition. Given a decomposition, they determine to which processor(s) a new item should be assigned. Currently, they work in conjunction with only the [RCB](#), [RIB](#), and [HSFC](#) algorithms.

[Zoltan_LB_Point_PP_Assign](#)
[Zoltan_LB_Box_PP_Assign](#)

For [backward compatibility](#) with previous versions of Zoltan, the following functions are also maintained. These functions are applicable only when the number of parts to be generated is equal to the number of processors on which the parts are computed. That is, these functions assume "parts" and "processors" are synonymous.

[Zoltan_LB_Point_Assign](#)
[Zoltan_LB_Box_Assign](#)

C:

int **Zoltan_LB_Point_PP_Assign** (
 struct **Zoltan_Struct** * *zz*,
 double * *coords*,
 int * *proc*,
 int * *part*);

FORTTRAN:

FUNCTION **Zoltan_LB_Point_PP_Assign**(*zz*, *coords*, *proc*, *part*)
INTEGER(Zoltan_INT) :: Zoltan_LB_Point_PP_Assign
TYPE(Zoltan_Struct), INTENT(IN) :: *zz*
REAL(Zoltan_DOUBLE), DIMENSION(*), INTENT(IN) :: *coords*
INTEGER(Zoltan_INT), INTENT(OUT) :: *proc*
INTEGER(Zoltan_INT), INTENT(OUT) :: *part*

C++:

int **Zoltan::LB_Point_PP_Assign** (
 double * const *coords*,
 int & *proc*,
 int & *part*);

Zoltan_LB_Point_PP_Assign is used to determine to which processor and parts a new point should be assigned. It is applicable only to geometrically generated decompositions ([RCB](#), [RIB](#), and [HSFC](#)). If the parameter **KEEP_CUTS** is set to TRUE, then the sequence of cuts that define the decomposition is saved. Given a new geometric point, the processor and parts which own it can be determined.

Arguments:

- zz*

Pointer to the Zoltan structure created by [Zoltan_Create](#).
- coords*

The (x,y) or (x,y,z) coordinates of the point being assigned.
- proc*

Upon return, the ID of the processor to which the point should belong.
- part*

Upon return, the ID of the parts to which the point should belong.

Returned Value:

- int

[Error code](#).

C:

int **Zoltan_LB_Box_PP_Assign** (

```

struct Zoltan_Struct * zz,
double xmin,
double ymin,
double zmin,
double xmax,
double ymax,
double zmax,
int *procs,
int *numprocs,
int *parts,
int *numparts);

```

FORTTRAN: **FUNCTION Zoltan_LB_Box_PP_Assign**(*zz, xmin, ymin, zmin, xmax, ymax, zmax, procs,*
numprocs, parts, numparts)
INTEGER(Zoltan_INT) :: Zoltan_LB_Box_PP_Assign
TYPE(Zoltan_Struct), **INTENT**(IN) :: *zz*
REAL(Zoltan_DOUBLE), **INTENT**(IN) :: *xmin, ymin, zmin, xmax, ymax, zmax*
INTEGER(Zoltan_INT), **DIMENSION**(*), **INTENT**(OUT) :: *procs*
INTEGER(Zoltan_INT), **INTENT**(OUT) :: *numprocs*
INTEGER(Zoltan_INT), **DIMENSION**(*), **INTENT**(OUT) :: *parts*
INTEGER(Zoltan_INT), **INTENT**(OUT) :: *numparts*

C++: **int Zoltan::LB_Box_PP_Assign** (
 const double & *xmin*,
 const double & *ymin*,
 const double & *zmin*,
 const double & *xmax*,
 const double & *ymax*,
 const double & *zmax*,
 int * const *procs*,
 int & *numprocs*,
 int * const *parts*,
 int & *numparts*);

In many settings, it is useful to know which processors and parts might need to know about an extended geometric object. **Zoltan_LB_Box_PP_Assign** addresses this problem. Given a geometric decomposition of space (currently only [RCB](#), [RIB](#), and [HSFC](#) are supported), and given an axis-aligned box around the geometric object, **Zoltan_LB_Box_PP_Assign** determines which processors and parts own geometry that intersects the box. To use this routine, the parameter **KEEP_CUTS** must be set to TRUE when the decomposition is generated. This parameter will cause the sequence of geometric cuts to be saved, which is necessary for **Zoltan_LB_Box_PP_Assign** to do its job.

Note that if the parameter **REDUCE_DIMENSIONS** was set to TRUE and the geometry was determined to be degenerate when decomposition was calculated, then the calculation was performed on transformed coordinates. This means that **Zoltan_LB_Box_PP_Assign** must transform the supplied bounding box accordingly. The transformed vertices are bounded again, and the parts intersections are calculated in the transformed space on this new bounding box. The impact of this is that **Zoltan_LB_Box_PP_Assign** may return parts not actually intersecting the original bounding box, but it will not omit any parts intersecting the original bounding box.

Arguments:

- zz* Pointer to the Zoltan structure created by [Zoltan_Create](#).
- xmin, ymin, zmin* The coordinates of the lower extent of the bounding box around the object. If the geometry is two-dimensional, the *z* value is ignored.

<i>xmax, ymax, zmax</i>	The coordinates of the upper extent of the bounding box around the object. If the geometry is two-dimensional, the <i>z</i> value is ignored.
<i>procs</i>	The list of processors intersecting the box are returned starting at this address. Note that <i>it is the responsibility of the calling routine to ensure that there is sufficient space for the return list</i> .
<i>numprocs</i>	Upon return, this value contains the number of processors that intersect the box (i.e. the number of entries placed in the <i>procs</i> list).
<i>parts</i>	The list of parts intersecting the box are returned starting at this address. Note that <i>it is the responsibility of the calling routine to ensure that there is sufficient space for the return list</i> .
<i>numparts</i>	Upon return, this value contains the number of parts that intersect the box (i.e. the number of entries placed in the <i>parts</i> list).

Returned Value:

int [Error code](#).

```
C:      int Zoltan_LB_Point_Assign (  
        struct Zoltan_Struct * zz,  
        double * coords,  
        int * proc);  
  
FORTRAN:  FUNCTION Zoltan_LB_Point_Assign(zz, coords, proc)  
           INTEGER(Zoltan_INT) :: Zoltan_LB_Point_Assign  
           TYPE(Zoltan_Struct), INTENT(IN) :: zz  
           REAL(Zoltan_DOUBLE), DIMENSION(*), INTENT(IN) :: coords  
           INTEGER(Zoltan_INT), INTENT(OUT) :: proc
```

Zoltan_LB_Point_Assign is is a wrapper around [Zoltan_LB_Point_PP_Assign](#) that excludes the parts assignment results. **Zoltan_LB_Point_Assign** assumes the number of parts is equal to the number of processors; thus, the parts assignment is equivalent to the processor assignment.

Arguments:

All arguments are analogous to those in [Zoltan_LB_Point_PP_Assign](#). Parts-assignment argument *part* is not included, as processor and parts numbers are considered to be the same in **Zoltan_LB_Point_Assign**.

Returned Value:

int [Error code](#).

```
C:      int Zoltan_LB_Box_Assign (  
        struct Zoltan_Struct * zz,  
        double xmin,  
        double ymin,  
        double zmin,  
        double xmax,  
        double ymax,  
        double zmax,  
        int *procs,  
        int *numprocs);  
  
FORTRAN:  FUNCTION Zoltan_LB_Box_Assign(zz, xmin, ymin, zmin, xmax, ymax, zmax, procs, numprocs)  
           INTEGER(Zoltan_INT) :: Zoltan_LB_Box_Assign  
           TYPE(Zoltan_Struct), INTENT(IN) :: zz
```

```
REAL(Zoltan_DOUBLE), INTENT(IN) :: xmin, ymin, zmin, xmax, ymax, zmax
INTEGER(Zoltan_INT), DIMENSION(*), INTENT(OUT) ::procs
INTEGER(Zoltan_INT), INTENT(OUT) :: numprocs
```

Zoltan_LB_Box_Assign is a wrapper around [Zoltan_LB_Box_PP_Assign](#) that excludes the parts assignment results. **Zoltan_LB_Box_Assign** assumes the number of parts is equal to the number of processors; thus, the parts assignment is equivalent to the processor assignment.

Arguments:

All arguments are analogous to those in [Zoltan_LB_Box_PP_Assign](#). Parts-assignment arguments *parts* and *numparts* are not included, as processor and parts numbers are considered to be the same in **Zoltan_LB_Box_Assign**.

Returned Value:

int [Error code](#).

Migration Functions

Zoltan's migration functions transfer object data to the processors in a new decomposition. Data to be transferred is specified through the import/export lists returned by [Zoltan_LB_Partition](#). Alternatively, users may specify their own import/export lists.

The migration functions can migrate objects based on their new part assignments and/or their new processor assignments. Behavior is determined by the [MIGRATE_ONLY_PROC_CHANGES](#) parameter.

If requested, Zoltan can automatically transfer an application's data between processors to realize a new decomposition. This functionality will be performed as part of the call to [Zoltan_LB_Partition](#) if the [AUTO_MIGRATE](#) parameter is set to TRUE (nonzero) via a call to [Zoltan_Set_Param](#). This approach is effective for when the data to be moved is relatively simple. For more complicated data movement, the application can leave [AUTO_MIGRATE](#) FALSE and call [Zoltan_Migrate](#) itself. In either case, [routines to pack and unpack object data](#) must be provided by the application. See the [Migration Examples](#) for examples with and without auto-migration.

The following functions are the migration interface functions. Their detailed descriptions can be found below.

[Zoltan_Invert_Lists](#) [Zoltan_Migrate](#)

The following functions are maintained for [backward compatibility](#) with previous versions of Zoltan. These functions are applicable only when the number of parts to be generated is equal to the number of processors on which the parts are computed. That is, these functions assume "parts" and "processors" are synonymous.

[Zoltan_Compute_Destinations](#) [Zoltan_Help_Migrate](#)

```

C:      int Zoltan_Invert_Lists (
        struct Zoltan_Struct *zz,
        int num_known,
        ZOLTAN\_ID\_PTR known_global_ids,
        ZOLTAN\_ID\_PTR known_local_ids,
        int *known_procs,
        int *known_to_part,
        int *num_found,
        ZOLTAN\_ID\_PTR *found_global_ids,
        ZOLTAN\_ID\_PTR *found_local_ids,
        int **found_procs,
        int **found_to_part);

FORTRAN: FUNCTION Zoltan_Invert_Lists(zz, num_known, known_global_ids, known_local_ids,
known_procs, known_to_part, num_found, found_global_ids, found_local_ids, found_procs,
found_to_part)
INTEGER(Zoltan_INT) :: Zoltan_Invert_Lists
TYPE(Zoltan_Struct),INTENT(IN) :: zz
INTEGER(Zoltan_INT), INTENT(IN) :: num_known
INTEGER(Zoltan_INT), INTENT(OUT) :: num_found
INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: known_global_ids, found_global_ids
INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: known_local_ids, found_local_ids

```

INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: known_procs, found_procs
INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: known_to_part, found_to_part

```
C++:
int Zoltan::Invert_Lists (
    const int & num_known,
    ZOLTAN\_ID\_PTR const known_global_ids,
    ZOLTAN\_ID\_PTR const known_local_ids,
    int * const known_procs,
    int * const known_to_part,
    int &num_found,
    ZOLTAN\_ID\_PTR &found_global_ids,
    ZOLTAN\_ID\_PTR &found_local_ids,
    int * &found_procs,
    int * &found_to_part);
```

Zoltan_Invert_Lists computes inverse communication maps useful for migrating data. It can be used in two ways:

- Given a list of known off-processor objects to be received by a processor, compute a list of local objects to be sent by the processor to other processors; or
- Given a list of known local objects to be sent by a processor to other processors, compute a list of off-processor objects to be received by the processor.

For example, if each processor knows which objects it will import from other processors, **Zoltan_Invert_Lists** computes the list of objects each processor needs to export to other processors. If, instead, each processor knows which objects it will export to other processors, **Zoltan_Invert_Lists** computes the list of objects each processor will import from other processors. The computed lists are allocated in Zoltan; they should not be allocated by the application before calling **Zoltan_Invert_Lists**. These lists can be freed through a call to [Zoltan_LB_Free_Part](#).

Arguments:

<i>zz</i>	Pointer to the Zoltan structure, created by Zoltan_Create , to be used in this invocation of the migration routine.
<i>num_known</i>	The number of known objects to be received (sent) by this processor.
<i>known_global_ids</i>	An array of <i>num_known</i> global IDs of known objects to be received (sent) by this processor. (size = <i>num_known</i> * NUM_GID_ENTRIES)
<i>known_local_ids</i>	An array of <i>num_known</i> local IDs of known objects to be received (sent) by this processor. (size = <i>num_known</i> * NUM_LID_ENTRIES)
<i>known_procs</i>	An array of size <i>num_known</i> listing the processor IDs of the processors that the known objects will be received from (sent to).
<i>known_to_part</i>	An array of size <i>num_known</i> listing the part numbers of the parts that the known objects will be assigned to.
<i>num_found</i>	Upon return, the number of objects that must be sent to (received from) other processors.
<i>found_global_ids</i>	Upon return, an array of <i>num_found</i> global IDs of objects to be sent (received) by this processor. (size = <i>num_found</i> * NUM_GID_ENTRIES)
<i>found_local_ids</i>	Upon return, an array of <i>num_found</i> local IDs of objects to be sent (received) by this processor. (size = <i>num_found</i> * NUM_LID_ENTRIES)
<i>found_procs</i>	Upon return, an array of size <i>num_found</i> listing the processor IDs of processors that the found objects will be sent to (received from).
<i>found_to_part</i>	An array of size <i>num_found</i> listing the part numbers of the parts that the found objects will be assigned to.

Returned Value:

int [Error code](#).

Note that the number of global and local ID entries ([NUM_GID_ENTRIES](#) and [NUM_LID_ENTRIES](#)) should be set using [Zoltan_Set_Param](#) before calling **Zoltan_Invert_Lists**. All processors must have the same values for these two parameters.

C: int **Zoltan_Migrate** (
 struct **Zoltan_Struct** *zz,
 int num_import,
 [ZOLTAN_ID_PTR](#) import_global_ids,
 [ZOLTAN_ID_PTR](#) import_local_ids,
 int *import_procs,
 int *import_to_part,
 int num_export,
 [ZOLTAN_ID_PTR](#) export_global_ids,
 [ZOLTAN_ID_PTR](#) export_local_ids,
 int *export_procs,
 int *export_to_part);

FORTTRAN: FUNCTION **Zoltan_Migrate**(zz, num_import, import_global_ids, import_local_ids, import_procs,
 import_to_part, num_export, export_global_ids, export_local_ids, export_procs, export_to_part)
 INTEGER(Zoltan_INT) :: Zoltan_Migrate
 TYPE(Zoltan_Struct),INTENT(IN) :: zz
 INTEGER(Zoltan_INT), INTENT(IN) :: num_import, num_export
 INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_global_ids, export_global_ids
 INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_local_ids, export_local_ids
 INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_procs, export_procs
 INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_to_part, export_to_part

C++: int **Zoltan::Migrate** (
 const int & num_import,
 [ZOLTAN_ID_PTR](#) const import_global_ids,
 [ZOLTAN_ID_PTR](#) const import_local_ids,
 int * const import_procs,
 int * const import_to_part,
 const int & num_export,
 [ZOLTAN_ID_PTR](#) const export_global_ids,
 [ZOLTAN_ID_PTR](#) const export_local_ids,
 int * const export_procs,
 int * const export_to_part);

Zoltan_Migrate takes lists of objects to be sent to other processors, along with the destinations of those objects, and performs the operations necessary to send the data associated with those objects to their destinations. **Zoltan_Migrate** performs the following operations using the application-registered functions:

- Call [ZOLTAN_PRE_MIGRATE_PP_FN_TYPE](#) (if registered)
- For each export object, call [ZOLTAN_OBJ_SIZE_FN_TYPE](#) to get object sizes.
- For each export object, call [ZOLTAN_PACK_OBJ_FN_TYPE](#) to load communication buffers.
- Communicate buffers to destination processors.
- Call [ZOLTAN_MID_MIGRATE_PP_FN_TYPE](#) (if registered).
- For each imported object, call [ZOLTAN_UNPACK_OBJ_FN_TYPE](#) to move data from the buffer into the new processor's data structures.
- Call [ZOLTAN_POST_MIGRATE_PP_FN_TYPE](#) (if registered).

Either export lists or import lists must be specified for **Zoltan_Migrate**. Both export lists and import lists may be specified, but both are not required.

If export lists are provided, non-NULL values for input arguments *import_global_ids*, *import_local_ids*, *import_procs*, and *import_to_part* are optional. The values must be non-NULL only if no export lists are provided or if the import lists are used by the application callback functions [ZOLTAN_PRE MIGRATE PP FN](#), [ZOLTAN MID MIGRATE PP FN](#), and [ZOLTAN POST MIGRATE PP FN](#). If all processors pass NULL arguments for the import arrays, the value of *num_import* should be -1.

Similarly, if import lists are provided, non-NULL values for input arguments *export_global_ids*, *export_local_ids*, *export_procs*, and *export_to_part* are optional. The values must be non-NULL only if no import lists are provided or if the export lists are used by the application callback functions [ZOLTAN_PRE MIGRATE PP FN](#), [ZOLTAN MID MIGRATE PP FN](#), and [ZOLTAN POST MIGRATE PP FN](#). If all processors pass NULL arguments for the export arrays, the value of *num_export* should be -1. In this case, **Zoltan_Migrate** computes the export lists based on the import lists.

Arguments:

<i>zz</i>	Pointer to the Zoltan structure, created by Zoltan_Create , to be used in this invocation of the migration routine.
<i>num_import</i>	The number of objects to be imported to parts on this processor; these objects may be stored on other processors or may be moving to new parts within this processor. Use <i>num_import</i> =-1 if all processors do not specify import arrays.
<i>import_global_ids</i>	An array of <i>num_import</i> global IDs of objects to be imported to parts on this processor. (size = <i>num_import</i> * NUM_GID_ENTRIES). All processors may pass <i>import_global_ids</i> =NULL if export lists are provided and <i>import_global_ids</i> is not needed by callback functions ZOLTAN_PRE MIGRATE PP FN , ZOLTAN MID MIGRATE PP FN , and ZOLTAN POST MIGRATE PP FN .
<i>import_local_ids</i>	An array of <i>num_import</i> local IDs of objects to be imported to parts on this processor. (size = <i>num_import</i> * NUM_LID_ENTRIES) All processors may pass <i>import_local_ids</i> =NULL if export lists are provided and <i>import_local_ids</i> is not needed by callback functions ZOLTAN_PRE MIGRATE PP FN , ZOLTAN MID MIGRATE PP FN , and ZOLTAN POST MIGRATE PP FN .
<i>import_procs</i>	An array of size <i>num_import</i> listing the processor IDs of objects to be imported to parts on this processor (i.e., the source processors). All processors may pass <i>import_procs</i> =NULL if export lists are provided and <i>import_procs</i> is not needed by callback functions ZOLTAN_PRE MIGRATE PP FN , ZOLTAN MID MIGRATE PP FN , and ZOLTAN POST MIGRATE PP FN .
<i>import_to_part</i>	An array of size <i>num_import</i> listing the parts to which imported objects should be assigned. All processors may pass <i>import_to_part</i> =NULL if export lists are provided and <i>import_to_part</i> is not needed by callback functions ZOLTAN_PRE MIGRATE PP FN , ZOLTAN MID MIGRATE PP FN , and ZOLTAN POST MIGRATE PP FN .
<i>num_export</i>	The number of objects that were stored on this processor in the previous decomposition that are assigned to other processors or to different parts within this processor in the new decomposition. Use <i>num_export</i> =-1 if all processors do not specify export arrays.
<i>export_global_ids</i>	An array of <i>num_export</i> global IDs of objects to be exported to new parts. (size = <i>num_export</i> * NUM_GID_ENTRIES) All processors may pass <i>export_global_ids</i> =NULL if import lists are provided and <i>export_global_ids</i> is not needed by callback functions ZOLTAN_PRE MIGRATE PP FN , ZOLTAN MID MIGRATE PP FN , and ZOLTAN POST MIGRATE PP FN .
<i>export_local_ids</i>	An array of <i>num_export</i> local IDs of objects to be exported to new parts. (size = <i>num_export</i> * NUM_LID_ENTRIES)

All processors may pass *export_local_ids*=NULL if import lists are provided and *export_local_ids* is not needed by callback functions [ZOLTAN_PRE_MIGRATE_PP_FN](#), [ZOLTAN_MID_MIGRATE_PP_FN](#), and [ZOLTAN_POST_MIGRATE_PP_FN](#).

export_procs

An array of size *num_export* listing the processor IDs to which exported objects should be assigned (i.e., the destination processors).
All processors may pass *export_procs*=NULL if import lists are provided and *export_procs* is not needed by callback functions [ZOLTAN_PRE_MIGRATE_PP_FN](#), [ZOLTAN_MID_MIGRATE_PP_FN](#), and [ZOLTAN_POST_MIGRATE_PP_FN](#).

export_to_part

An array of size *num_export* listing the parts to which exported objects should be assigned. All processors may pass *export_to_part*=NULL if import lists are provided and *export_to_part* is not needed by callback functions [ZOLTAN_PRE_MIGRATE_PP_FN](#), [ZOLTAN_MID_MIGRATE_PP_FN](#), and [ZOLTAN_POST_MIGRATE_PP_FN](#).

Returned Value:

int [Error code](#).

Note that the number of global and local ID entries ([NUM_GID_ENTRIES](#) and [NUM_LID_ENTRIES](#)) should be set using [Zoltan_Set_Param](#) before calling **Zoltan_Migrate**. All processors must have the same values for these two parameters.

```
C:      int Zoltan_Compute_Destinations (
        struct Zoltan_Struct *zz,
        int num_known,
        ZOLTAN\_ID\_PTR known_global_ids,
        ZOLTAN\_ID\_PTR known_local_ids,
        int *known_procs,
        int *num_found,
        ZOLTAN\_ID\_PTR *found_global_ids,
        ZOLTAN\_ID\_PTR *found_local_ids,
        int **found_procs);

FORTRAN: FUNCTION Zoltan_Compute_Destinations(zz, num_known, known_global_ids,
        known_local_ids, known_procs, num_found, found_global_ids, found_local_ids, found_procs)
        INTEGER(Zoltan_INT) :: Zoltan_Compute_Destinations
        TYPE(Zoltan_Struct),INTENT(IN) :: zz
        INTEGER(Zoltan_INT), INTENT(IN) :: num_known
        INTEGER(Zoltan_INT), INTENT(OUT) :: num_found
        INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: known_global_ids, found_global_ids
        INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: known_local_ids, found_local_ids
        INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: known_procs, found_procs
```

Zoltan_Compute_Destinations is a wrapper around [Zoltan_Invert_Lists](#) that excludes part assignment arrays. It is maintained for backward compatibility with previous versions of Zoltan.

Zoltan_Compute_Destinations assumes the number of parts is equal to the number of processors. The computed lists are allocated in Zoltan; they should not be allocated by the application before calling **Zoltan_Compute_Destinations**. These lists can be freed through a call to [Zoltan_LB_Free_Data](#) or [Zoltan_LB_Free_Part](#).

Arguments:

All arguments are analogous to those in [Zoltan_Invert_Lists](#). Part-assignment arrays *known_to_part* and *found_to_part* are not included, as part and processor numbers are assumed to be the same in **Zoltan_Compute_Destinations**.

Returned Value:

int [Error code](#).

Note that the number of global and local ID entries ([NUM_GID_ENTRIES](#) and [NUM_LID_ENTRIES](#)) should be set using [Zoltan_Set_Param](#) before calling **Zoltan_Compute_Destinations**. All processors must have the same values for these two parameters.

```
C:      int Zoltan_Help_Migrate (  
        struct Zoltan_Struct *zz,  
        int num_import,  
        ZOLTAN\_ID\_PTR import_global_ids,  
        ZOLTAN\_ID\_PTR import_local_ids,  
        int *import_procs,  
        int num_export,  
        ZOLTAN\_ID\_PTR export_global_ids,  
        ZOLTAN\_ID\_PTR export_local_ids,  
        int *export_procs);  
  
FORTRAN:  FUNCTION Zoltan_Help_Migrate(zz, num_import, import_global_ids, import_local_ids,  
    import_procs, num_export, export_global_ids, export_local_ids, export_procs)  
    INTEGER(Zoltan_INT) :: Zoltan_Help_Migrate  
    TYPE(Zoltan_Struct),INTENT(IN) :: zz  
    INTEGER(Zoltan_INT), INTENT(IN) :: num_import, num_export  
    INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_global_ids, export_global_ids  
    INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_local_ids, export_local_ids  
    INTEGER(Zoltan_INT), POINTER, DIMENSION(:) :: import_procs, export_procs
```

Zoltan_Help_Migrate is a wrapper around [Zoltan_Migrate](#) that excludes part assignment arrays. It is maintained for backward compatibility with previous versions of Zoltan.

Zoltan_Help_Migrate assumes the number of parts is equal to the number of processors. It uses migration pre-, mid-, and post-processing routines [ZOLTAN_PRE_MIGRATE_FN_TYPE](#), [ZOLTAN_MID_MIGRATE_FN_TYPE](#), and [ZOLTAN_POST_MIGRATE_FN_TYPE](#), respectively, which also exclude part assignment arrays.

Arguments:

All arguments are analogous to those in [Zoltan_Migrate](#). Part-assignment arrays *import_to_part* and *export_to_part* are not included, as part and processor numbers are assumed to be the same in **Zoltan_Help_Migrate**.

Returned Value:

int [Error code](#).

Ordering Functions

Zoltan provides functions for ordering a set of objects, typically given as a graph (which may correspond to a sparse matrix). The following functions are the ordering interface functions in Zoltan. The first is the main function, and the others are accessor functions that should only be called after [Zoltan_Order](#).

- [Zoltan_Order](#)
- [Zoltan_Order_Get_Num_Blocks](#)
- [Zoltan_Order_Get_Block_Bounds](#)
- [Zoltan_Order_Get_Block_Size](#)
- [Zoltan_Order_Get_Block_Parent](#)
- [Zoltan_Order_Get_Num_Leaves](#)
- [Zoltan_Order_Get_Block_Leaves](#)

```

C:      int Zoltan_Order (
        struct Zoltan_Struct *zz,
        int num_gid_entries,
        int num_obj,
        ZOLTAN\_ID\_PTR global_ids,
        ZOLTAN\_ID\_PTR rank,
        )

FORTRAN: FUNCTION Zoltan_Order(zz, num_gid_entries, num_obj, global_ids, rank, iperm)
          INTEGER(Zoltan_INT) :: Zoltan_Order
          TYPE(Zoltan_Struct), INTENT(IN) :: zz
          INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries
          INTEGER(Zoltan_INT), INTENT(IN) :: num_obj
          INTEGER(Zoltan_INT) :: global_ids(*)
          INTEGER(Zoltan_INT) :: rank(*)

C++:    int Zoltan::Order (
        int num_gid_entries,
        int num_obj,
        ZOLTAN\_ID\_PTR global_ids,
        ZOLTAN\_ID\_PTR rank,
        )
    
```

Zoltan_Order invokes the ordering routine specified by the [ORDER_METHOD](#) parameter. Results of the ordering is returned in the array *rank* .

- *rank[i]* gives the rank of *global_ids[i]* in the computed ordering, which is a number between 0 and N-1 where N is the overall number of objects across all the processors. (Note: This will change in future versions. A permuted set of GIDs will be returned instead.)
- The arrays *global_ids*, *rank*, should all be allocated by the application before **Zoltan_Order** is called. Each array must have space for (at least) *num_obj* elements, where *num_obj* is the number of objects the user want to know informations about.

Arguments:

zz Pointer to the Zoltan structure, created by [Zoltan_Create](#), to be used in this invocation of the

load-balancing routine.

<i>num_gid_entries</i>	Input: the number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_obj</i>	Number of objects for which we want to know the ordering. Objects may be non-local.
<i>global_ids</i>	An array of global IDs of objects for which we want to know the ordering on this processor. Size of this array must be <i>num_obj</i> . Objects may be non-local. Objects IDs may be repeated on several processor.
<i>rank</i>	Upon return, an array of length <i>num_obj</i> containing the rank of each object in the computed ordering. When <i>rank[i] = j</i> , that means that the object corresponding to <i>global_ids[i]</i> is the <i>j</i> th object in the ordering. (This array corresponds directly to the <i>perm</i> array in METIS and the <i>order</i> array in ParMETIS.) Note that the rank may refer to either a local or a global ordering, depending on ORDER_TYPE . Memory for this array must have been allocated before Zoltan_Order is called.

Returned Value:

int [Error code](#).

Accessors

Zoltan primarily supports nested dissection orderings, which are typically used to reduce fill in direct solvers. For this use case, it is important to get additional information (the separators). The accessor functions below define the Zoltan interface. Note that these functions should be called after [Zoltan_Order](#).

C:	int Zoltan_Order_Get_Num_Blocks (struct Zoltan_Struct *zz)
C++:	int Zoltan::Order_Get_Num_Blocks ()

Zoltan_Order_Get_Num_Blocks returns the number of subdomains (or column blocks) computed during the ordering. For a Nested Dissection based ordering method, it corresponds to the sum of the separators and the subgraphs of the lowest level. For example, after one bisection the number of blocks is three.

Arguments:

zz	Pointer to the Zoltan structure, created by Zoltan_Create , which has been used in the ordering routine.
----	--------------------------------------------------------------------------------------------------------------------------

Returned Value:

int The number of *blocks* (subdomains).

C:	int Zoltan_Order_Get_Block_Bounds (struct Zoltan_Struct *zz, int <i>block_id</i> , int * <i>first</i> , int * <i>last</i>)
C++:	int Zoltan::Order_Get_Block_Bounds (int <i>block_id</i> , int & <i>first</i> , int & <i>last</i>)

Zoltan_Order_Get_Block_Bounds gives the boundaries of the given block. The *first* and *last* parameters contain

upon return the indices of the begin and the end of the block. These indices are from the global continuous numbering between 1 and N-1 of the N distributed objects.

Arguments:

<i>zz</i>	Pointer to the Zoltan structure, created by Zoltan_Create , which has been used in the ordering routine.
<i>block_id</i>	The number of the block we want informations on.
<i>first</i>	Upon return, pointer to the value of the begining of the block.
<i>last</i>	Upon return, pointer to the value of the end of the block.

Returned Value:

int	Error code .
-----	------------------------------

C:	int Zoltan_Order_Get_Block_Size (struct Zoltan_Struct * <i>zz</i> , int <i>block_id</i>)
C++:	int Zoltan::Order_Get_Block_Size (int <i>block_id</i>)

Zoltan_Order_Get_Block_Size returns the number of objects in the block *block_id*.

Arguments:

<i>zz</i>	Pointer to the Zoltan structure, created by Zoltan_Create , which has been used in the ordering routine.
<i>block_id</i>	The indice of the block we want to know the size.

Returned Value:

int	The number of objects in this given <i>block</i> .
-----	----------------------------------------------------

C:	int Zoltan_Order_Get_Block_Parent (struct Zoltan_Struct * <i>zz</i> , int <i>block_id</i>)
C++:	int Zoltan::Order_Get_Block_Parent (int <i>block_id</i>)

Zoltan_Order_Get_Block_Parent returns the number (id) of the parent block in the elimination tree. The value is -1 for the root of the tree.

Arguments:

<i>zz</i>	Pointer to the Zoltan structure, created by Zoltan_Create , which has been used in the ordering routine.
<i>block_id</i>	The indice of the block we want to know the size.

Returned Value:

int	The number (id) of the parent block in the elimination tree. The value is -1 for the root of the tree.
-----	--------------------------------------------------------------------------------------------------------

```
C:          int Zoltan_Order_Get_Num_Leaves (  
            struct Zoltan_Struct *zz)  
  
C++:       int Zoltan::Order_Get_Num_Leaves ()
```

Zoltan_Order_Get_Num_Leaves returns the number of leaves in the elimination tree.

Arguments:

zz Pointer to the Zoltan structure, created by [Zoltan_Create](#), which has been used in the ordering routine.

Returned Value:

int The number of leaves in the elimination tree.

```
C:          void Zoltan_Order_Get_Block_Leaves (  
            struct Zoltan_Struct *zz,  
            int *leaves)  
  
C++:       void Zoltan::Order_Get_Block_Leaves (  
            struct Zoltan_Struct *zz,  
            int *leaves)
```

Zoltan_Order_Get_Block_Leaves get the indices of the blocks that are leaves in the elimination tree.

Arguments:

zz Pointer to the Zoltan structure, created by [Zoltan_Create](#), which has been used in the ordering routine.

leaves Array of indices of the blocks that are leaves in the elimination. The last element of this array is -1. The array must be of size *number of leave + 1* and must be allocated by the user before the call.

Coloring Functions

Zoltan provides limited capability for coloring a set of objects, typically given as a graph. In graph coloring, each vertex is assigned an integer label such that no two adjacent vertices have the same label. The following functions are the coloring interface functions in the Zoltan library; their descriptions are included below.

[Zoltan_Color](#)

C:

int **Zoltan_Color** (
 struct **Zoltan_Struct** *zz,
 int num_gid_entries,
 int num_obj,
 [ZOLTAN_ID_PTR](#) global_ids,
 int *color_exp);

FORTTRAN:

Not yet available.

C++:

int **Zoltan::Color** (
 int &num_gid_entries,
 const int &num_obj,
 [ZOLTAN_ID_PTR](#) global_ids,
 int *color_exp);

Zoltan_Color invokes the coloring routine and the assigned colors of each object are returned in the array *color_exp*. *color_exp[i]* gives the color of *global_ids[i]* in the computed coloring. The arrays *global_ids* and *color_exp* should all be allocated by the application before **Zoltan_Color** is called. *global_ids* must contain the global ids of the elements for which the current processor wants coloring informations.

Arguments:

<i>zz</i>	Pointer to the Zoltan structure, created by Zoltan_Create , to be used in this invocation of the load-balancing routine.
<i>num_gid_entries</i>	Input: the number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_obj</i>	Number of objects for which we want to know the color on this processor. Objects may be non-local or duplicated.
<i>global_ids</i>	An array of global IDs of objects for which we want to know the color on this processor. Size of this array must be <i>num_obj</i> . Objects may be non-local. Objects IDs may be repeated on several processor.
<i>color_exp</i>	Upon return, an array of length <i>num_obj</i> containing the colors of objects. That is, <i>color_exp[i]</i> gives the color of <i>global_ids[i]</i> in the computed coloring. By default, colors are positive integers starting at one. Memory for this array must have been allocated before Zoltan_Color is called.

Returned Value:

int	Error code .
-----	------------------------------

Application-Registered Query Functions

Zoltan gets information about a processor's objects through calls to query functions. These functions must be provided by the application. They are "registered" with Zoltan; that is, a pointer to the function is passed to Zoltan, which can then call that function when its information is needed.

Query functions return information about only on-processor data. They can be called by Zoltan with individual objects or lists of objects. Each processor may call a given query function zero, one or more than one time. Thus, most query functions should NOT contain interprocessor communication, as such communication can cause processors to hang while waiting for messages that were never sent. (The only exceptions to this rule are certain [migration query functions](#).)

Two categories of query functions are used by the library:

[General Zoltan Query Functions](#)
[Migration Query Functions](#)

In each category, a variety of query functions can be registered by the user. The query functions have a function type, describing their purpose. Functions can be registered with a Zoltan structure in two ways: through calls to [Zoltan_Set_Fn](#) or through calls to query-function-specific functions [Zoltan_Set <zoltan_fn_type> Fn](#). When a function is registered through a call to [Zoltan_Set_Fn](#), its function type is passed in the *fn_type* argument. When [Zoltan_Set <zoltan_fn_type> Fn](#) is used to register functions, the type of the function is implicit in the *fn_ptr* argument. Each function description below includes both its function type and function prototype.

Query functions that return information about data objects owned by a processor come in two forms: list-based functions that return information about a list of objects, and iterator functions that return information about a single object. Users can provide either version of the query function; they need not provide both. Zoltan calls the list-based functions with the IDs of all objects needed; this approach often provides faster performance as it eliminates the overhead of multiple function calls. List-based functions have the word "MULTI" in their function-type name. If, instead, the application provides iterator functions, Zoltan calls the iterator function once for each object whose data is needed. This approach, while slower, allows Zoltan to use less memory for some data.

Some algorithms in Zoltan require that certain query functions be registered by the application; for example, geometric partitioning algorithms such as Recursive Coordinate Bisection (RCB) require that either a [ZOLTAN_GEOM_FN](#) or a [ZOLTAN_GEOM_MULTI_FN](#) be registered. When a default value is specified below, the query function type is optional; if a function of that type is not registered, the default values are used. Details of which query functions are required by particular algorithms are included in the [Algorithms](#) section.

Many of the functions have both global and local object identifiers (IDs) in their argument lists. The global IDs provided by the application must be unique across all processors; they are used for identification within Zoltan. The local IDs are not used by Zoltan; they are provided for the convenience of the application and can be anything the application desires. The local IDs can be used by application query routines to enable direct access to application data. For example, the object with global ID "3295" may be stored by the application in location "15" of an array in the processor's local memory. Both global ID "3295" and local ID "15" can be used by the application to describe the object. Then, rather than searching the array for global ID "3295," the application query routines can subsequently use the local ID to index directly into the local storage array. See [Data Types for Object IDs](#) for a description of global and local IDs. All of the functions have, as their first argument, a pointer to data that is passed to Zoltan through [Zoltan_Set_Fn](#) or [Zoltan_Set <zoltan_fn_type> Fn](#). This data is not used by Zoltan. A different set of data can be supplied for each registered function. For example, if the local ID is an index into an array of data structures, then the data pointer might point to the head of the data structure array.

As their last argument, all functions have an [error code](#) that should be set and returned by the registered function.

If you are calling the Zoltan library from a C++ application, you may set the query function to be any class static function or any function defined outside of a class definition. However, it is possible you will wish to set the query function to be an object method. In that case, you should write a query function that takes a pointer to the object as its *data* field. The query function can then call the object method.

[\[Table of Contents\]](#) | [\[Next: Load-Balancing Query Functions\]](#) | [\[Previous: Coloring Functions\]](#) | [\[Privacy and Security\]](#)

General Zoltan Query Functions

The following registered functions are used by various Zoltan algorithms in the Zoltan library. No single algorithm uses all the query functions; the [algorithm descriptions](#) indicate which query functions are required by individual algorithms. These query functions should NOT contain interprocessor communication.

[Object ID Functions](#)

[ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)
[ZOLTAN_FIRST_OBJ_FN](#) (deprecated)
[ZOLTAN_NEXT_OBJ_FN](#) (deprecated)
[ZOLTAN_PART_MULTI_FN](#) or [ZOLTAN_PART_FN](#)

[Geometry-Based Functions](#)

[ZOLTAN_NUM_GEOM_FN](#)
[ZOLTAN_GEOM_MULTI_FN](#) or [ZOLTAN_GEOM_FN](#)

[Graph-Based Functions](#)

[ZOLTAN_NUM_EDGES_MULTI_FN](#) or [ZOLTAN_NUM_EDGES_FN](#)
[ZOLTAN_EDGE_LIST_MULTI_FN](#) or [ZOLTAN_EDGE_LIST_FN](#)

[Hypergraph-Based Functions](#)

[ZOLTAN_HG_SIZE_CS_FN](#)
[ZOLTAN_HG_CS_FN](#)
[ZOLTAN_HG_SIZE_EDGE_WTS_FN](#)
[ZOLTAN_HG_EDGE_WTS_FN](#)
[ZOLTAN_NUM_FIXED_OBJ_FN](#)
[ZOLTAN_FIXED_OBJ_LIST_FN](#)

[Tree-Based Functions](#)

[ZOLTAN_NUM_COARSE_OBJ_FN](#)
[ZOLTAN_COARSE_OBJ_LIST_FN](#)
[ZOLTAN_FIRST_COARSE_OBJ_FN](#)
[ZOLTAN_NEXT_COARSE_OBJ_FN](#)
[ZOLTAN_NUM_CHILD_FN](#)
[ZOLTAN_CHILD_LIST_FN](#)
[ZOLTAN_CHILD_WEIGHT_FN](#)

[Hierarchical Partitioning Functions](#)

[ZOLTAN_HIER_NUM_LEVELS_FN](#)
[ZOLTAN_HIER_PART_FN](#)
[ZOLTAN_HIER_METHOD_FN](#)

Object ID Functions

C and C++: typedef int **ZOLTAN_NUM_OBJ_FN** (void **data*, int **ierr*);

FORTRAN: FUNCTION *Get_Num_Obj*(*data*, *ierr*)
 INTEGER(Zoltan_INT) :: Get_Num_Obj
 <*type-data*>, INTENT(IN) :: *data*
 INTEGER(Zoltan_INT), INTENT(OUT) :: *ierr*

 <*type-data*> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),
 DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_*x*)
 where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_NUM_OBJ_FN** query function returns the number of objects that are currently assigned to the processor.

Function Type: **ZOLTAN_NUM_OBJ_FN_TYPE**

Arguments:

data Pointer to user-defined data.

ierr Error code to be set by function.

Returned Value:

 int The number of objects that are assigned to the processor.

C and C++: typedef void **ZOLTAN_OBJ_LIST_FN** (void **data*, int *num_gid_entries*, int *num_lid_entries*,
 [ZOLTAN_ID_PTR](#) *global_ids*, [ZOLTAN_ID_PTR](#) *local_ids*, int *wgt_dim*, float **obj_wgts*,
 int **ierr*);

FORTRAN: SUBROUTINE *Get_Obj_List*(*data*, *num_gid_entries*, *num_lid_entries*, *global_ids*, *local_ids*,
 wgt_dim, *obj_wgts*, *ierr*)
 <*type-data*>, INTENT(IN) :: *data*
 INTEGER(Zoltan_INT), INTENT(IN) :: *num_gid_entries*, *num_lid_entries*
 INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: *global_ids*
 INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: *local_ids*
 INTEGER(Zoltan_INT), INTENT(IN) :: *wgt_dim*
 REAL(Zoltan_FLOAT), INTENT(OUT), DIMENSION(*) :: *obj_wgts*
 INTEGER(Zoltan_INT), INTENT(OUT) :: *ierr*

 <*type-data*> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),
 DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_*x*)
 where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_OBJ_LIST_FN** query function fills two (three if weights are used) arrays with information about the objects currently assigned to the processor. Both arrays are allocated (and subsequently freed) by Zoltan; their size is determined by a call to a [ZOLTAN_NUM_OBJ_FN](#) query function to get the array size. For many algorithms, either a **ZOLTAN_OBJ_LIST_FN** query function or a [ZOLTAN_FIRST_OBJ_FN/ZOLTAN_NEXT_OBJ_FN](#) query-function pair must be registered; however, both query options need not be provided. The **ZOLTAN_OBJ_LIST_FN** is preferred for efficiency.

Function Type: **ZOLTAN_OBJ_LIST_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>global_ids</i>	Upon return, an array of unique global IDs for all objects assigned to the processor.
<i>local_ids</i>	Upon return, an array of local IDs, the meaning of which can be determined by the application, for all objects assigned to the processor. (Optional.)
<i>wgt_dim</i>	The number of weights associated with an object (typically 1), or 0 if weights are not requested. This value is set through the parameter OBJ_WEIGHT_DIM .
<i>obj_wgts</i>	Upon return, an array of object weights. Weights for object <i>i</i> are stored in <i>obj_wgts[(i-1)*wgt_dim:i*wgt_dim-1]</i> . If <i>wgt_dim</i> =0, the return value of <i>obj_wgts</i> is undefined and may be NULL.
<i>ierr</i>	Error code to be set by function.

C and C++:	<pre>typedef int ZOLTAN_FIRST_OBJ_FN (void *data, int num_gid_entries, int num_lid_entries, ZOLTAN_ID_PTR first_global_id, ZOLTAN_ID_PTR first_local_id, int wgt_dim, float *first_obj_wgt, int *ierr);</pre>
FORTRAN:	<pre>FUNCTION Get_First_Obj(data, num_gid_entries, num_lid_entries, first_global_id, first_local_id, wgt_dim, first_obj_wgt, ierr) INTEGER(Zoltan_INT) :: Get_First_Obj <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: first_global_id INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: first_local_id INTEGER(Zoltan_INT), INTENT(IN) :: wgt_dim REAL(Zoltan_FLOAT), INTENT(OUT), DIMENSION(*) :: first_obj_wgt INTEGER(Zoltan_INT), INTENT(OUT) :: ierr</pre> <p><type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where <i>x</i> is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</p>

(Deprecated) A **ZOLTAN_FIRST_OBJ_FN** query function initializes an iteration over objects assigned to the processor. It returns the global and local IDs of the first object on the processor. Subsequent calls to a [ZOLTAN_NEXT_OBJ_FN](#) query function iterate over and return other objects assigned to the processor. For many algorithms, either a [ZOLTAN_OBJ_LIST_FN](#) query function or a **ZOLTAN_FIRST_OBJ_FN**/[ZOLTAN_NEXT_OBJ_FN](#) query-function pair must be registered; however, both query options need not be provided. The [ZOLTAN_OBJ_LIST_FN](#) is preferred for efficiency.

Function Type: **ZOLTAN_FIRST_OBJ_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local

ids are not used.)

<i>first_global_id</i>	The returned value of the global ID for the first object; the value is ignored if there are no objects.
<i>first_local_id</i>	The returned value of the local ID for the first object; the value is ignored if there are no objects.
<i>wgt_dim</i>	The number of weights associated with an object (typically 1), or 0 if weights are not requested. This value is set through the parameter OBJ_WEIGHT_DIM .
<i>first_obj_wgt</i>	Upon return, the first object's weights; an array of length <i>wgt_dim</i> . Undefined if <i>wgt_dim</i> =0.
<i>ierr</i>	Error code to be set by function.

Returned Value:

1	If <i>first_global_id</i> and <i>first_local_id</i> contain valid IDs of the first object.
0	If no objects are available.

C and C++:	<pre>typedef int ZOLTAN_NEXT_OBJ_FN (void * data, int num_gid_entries, int num_lid_entries, ZOLTAN_ID_PTR global_id, ZOLTAN_ID_PTR local_id, ZOLTAN_ID_PTR next_global_id, ZOLTAN_ID_PTR next_local_id, int wgt_dim, float *next_obj_wgt, int *ierr);</pre>
FORTTRAN:	<pre>FUNCTION <i>Get_Next_Obj</i>(data, num_gid_entries, num_lid_entries, global_id, local_id, next_global_id, next_local_id, wgt_dim, next_obj_wgt, ierr) INTEGER(Zoltan_INT) :: Get_Next_Obj <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_id INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: local_id INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: next_global_id INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: next_local_id INTEGER(Zoltan_INT), INTENT(IN) :: wgt_dim REAL(Zoltan_FLOAT), INTENT(OUT), DIMENSION(*) :: next_obj_wgt INTEGER(Zoltan_INT), INTENT(OUT) :: ierr <type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where x is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</pre>

(Deprecated) A **ZOLTAN_NEXT_OBJ_FN** query function is an iterator function which, when given an object assigned to the processor, returns the next object assigned to the processor. The first object of the iteration is provided by a [ZOLTAN_FIRST_OBJ_FN](#) query function. For many algorithms, either a [ZOLTAN_OBJ_LIST_FN](#) query function or a [ZOLTAN_FIRST_OBJ_FN/ZOLTAN_NEXT_OBJ_FN](#) query-function pair must be registered; however, both query options need not be provided. The [ZOLTAN_OBJ_LIST_FN](#) is preferred for efficiency.

Function Type: **ZOLTAN_NEXT_OBJ_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>global_id</i>	The global ID of the previous object.

<i>local_id</i>	The local ID of the previous object.
<i>next_global_id</i>	The returned value of the global ID for the next object; the value is ignored if there are no more objects.
<i>next_local_id</i>	The returned value of the local ID for the next object; the value is ignored if there are no more objects.
<i>wgt_dim</i>	The number of weights associated with an object (typically 1), or 0 if weights are not requested. This value is set through the parameter OBJ_WEIGHT_DIM .
<i>next_obj_wgt</i>	Upon return, the next object's weights; an array of length <i>wgt_dim</i> . Undefined if <i>wgt_dim</i> =0.
<i>ierr</i>	Error code to be set by function.
Returned Value:	
1	If <i>next_global_id</i> and <i>next_local_id</i> contain valid IDs of the next object.
0	If no more objects are available.

C and C++:	typedef void ZOLTAN_PART_MULTI_FN (void *data, int num_gid_entries, int num_lid_entries, int num_obj, ZOLTAN_ID_PTR global_ids, ZOLTAN_ID_PTR local_ids, int *parts, int *ierr);
FORTRAN:	<pre>SUBROUTINE <i>Get_Part_Multi</i>(data, num_gid_entries, num_lid_entries, num_obj, global_ids, local_ids, ierr) <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries, num_obj INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_ids INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: local_ids INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: parts INTEGER(Zoltan_INT), INTENT(OUT) :: ierr</pre> <p><type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where <i>x</i> is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</p>

A **ZOLTAN_PART_MULTI_FN** query function returns a list of parts to which given objects are currently assigned. If a **ZOLTAN_PART_MULTI_FN** or [ZOLTAN_PART_FN](#) is not registered, Zoltan assumes the part numbers are the processor number of the owning processor. Valid part numbers are non-negative integers. This function is used when parameter [REMAP](#)=1 and for certain methods with parameter [LB_APPROACH](#)=*Repartition*.

Function Type: **ZOLTAN_PART_MULTI_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>num_obj</i>	The number of object IDs in arrays <i>global_ids</i> and <i>local_ids</i> .
<i>global_ids</i>	The global IDs of the objects for which the part numbers should be returned.
<i>local_ids</i>	The local IDs of the objects for which the part numbers should be returned. (Optional.)
<i>parts</i>	Upon return, an array of part numbers corresponding to the global and local IDs.
<i>ierr</i>	Error code to be set by function.

C and C++:

typedef int **ZOLTAN_PART_FN** (void **data*, int *num_gid_entries*, int *num_lid_entries*,
[ZOLTAN_ID_PTR](#) *global_id*, [ZOLTAN_ID_PTR](#) *local_id*, int **ierr*);

FORTRAN:

FUNCTION *Get_Part*(*data*, *num_gid_entries*, *num_lid_entries*, *global_id*, *local_id*, *ierr*)
 INTEGER(Zoltan_INT) :: Get_Part
 <*type-data*>, INTENT(IN) :: *data*
 INTEGER(Zoltan_INT), INTENT(IN) :: *num_gid_entries*, *num_lid_entries*
 INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: *global_id*
 INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: *local_id*
 INTEGER(Zoltan_INT), INTENT(OUT) :: *ierr*

<*type-data*> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),
DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_*x*)
where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_PART_FN** query function returns the part to which a given object is currently assigned. If a **ZOLTAN_PART_FN** or [ZOLTAN_PART_MULTI_FN](#) is not registered, Zoltan assumes the part numbers are the processor number of the owning processor. Valid part numbers are non-negative integers. This function is used when parameter [REMAP](#)=1 and for certain methods with parameter [LB_APPROACH](#)=*Repartition*.

Function Type:	ZOLTAN_PART_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>global_id</i>	The global ID of the object for which the part number should be returned.
<i>local_id</i>	The local ID of the object for which the part number should be returned.
<i>ierr</i>	Error code to be set by function.
Returned Value:	
int	The part number for the object identified by <i>global_id</i> and <i>local_id</i> .

Geometry-based Functions

C and C++:

typedef int **ZOLTAN_NUM_GEOM_FN** (void **data*, int **ierr*);

FORTRAN:

FUNCTION *Get_Num_Geom*(*data*, *ierr*)
 INTEGER(Zoltan_INT) :: Get_Num_Geom
 <*type-data*>, INTENT(IN) :: *data*
 INTEGER(Zoltan_INT), INTENT(OUT) :: *ierr*

<*type-data*> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),
DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_*x*)
where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_NUM_GEOM_FN** query function returns the number of values needed to express the geometry of an object. For example, for a two-dimensional mesh-based application, (x,y) coordinates are needed to describe an object's geometry; thus the **ZOLTAN_NUM_GEOM_FN** query function should return the value of two. For a similar three-dimensional application, the return value should be three.

Function Type: **ZOLTAN_NUM_GEOM_FN_TYPE**

Arguments:

- data* Pointer to user-defined data.
- ierr* Error code to be set by function.

Returned Value:

- int The number of values needed to express the geometry of an object.

```
C and C++:       typedef void ZOLTAN_GEOM_MULTI_FN (void *data, int num_gid_entries,
int num_lid_entries, int num_obj, ZOLTAN\_ID\_PTR global_ids, ZOLTAN\_ID\_PTR local_ids,
int num_dim, double *geom_vec, int *ierr);

FORTRAN:       SUBROUTINE Get_Geom_Multi(data, num_gid_entries, num_lid_entries, num_obj, global_ids,
local_ids, num_dim, geom_vec, ierr)
<type-data>, INTENT(IN) :: data
INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries
INTEGER(Zoltan_INT), INTENT(IN) :: num_obj, num_dim
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_ids
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: local_ids
REAL(Zoltan_DOUBLE), INTENT(OUT), DIMENSION(*) :: geom_vec
INTEGER(Zoltan_INT), INTENT(OUT) :: ierr

<type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),
DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x)
where x is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.
```

A **ZOLTAN_GEOM_MULTI_FN** query function returns a vector of geometry values for a list of given objects. The geometry vector is allocated by Zoltan to be of size *num_obj * num_dim*; its format is described below.

Function Type: **ZOLTAN_GEOM_MULTI_FN_TYPE**

Arguments:

- data* Pointer to user-defined data.
- num_gid_entries* The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter [NUM_GID_ENTRIES](#).
- num_lid_entries* The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter [NUM_LID_ENTRIES](#). (It should be zero if local ids are not used.)
- num_obj* The number of object IDs in arrays *global_ids* and *local_ids*.
- global_ids* Array of global IDs of objects whose geometry values should be returned.
- local_ids* Array of local IDs of objects whose geometry values should be returned. (Optional.)
- num_dim* Number of coordinate entries per object (typically 1, 2, or 3).
- geom_vec* Upon return, an array containing geometry values. For object *i* (specified by *global_ids[i*num_gid_entries]* and *local_ids[i*num_lid_entries]*, *i=0,1,...,num_obj-1*), coordinate values should be stored in *geom_vec[i*num_dim:(i+1)*num_dim-1]*.

<i>ierr</i>	Error code to be set by function.
<hr/>	
C and C++:	typedef void ZOLTAN_GEOM_FN (void * <i>data</i> , int <i>num_gid_entries</i> , int <i>num_lid_entries</i> , ZOLTAN_ID_PTR <i>global_id</i> , ZOLTAN_ID_PTR <i>local_id</i> , double * <i>geom_vec</i> , int * <i>ierr</i>);
FORTRAN:	SUBROUTINE <i>Get_Geom</i> (<i>data</i> , <i>num_gid_entries</i> , <i>num_lid_entries</i> , <i>global_id</i> , <i>local_id</i> , <i>geom_vec</i> , <i>ierr</i>) < <i>type-data</i> >, INTENT(IN) :: <i>data</i> INTEGER(Zoltan_INT), INTENT(IN) :: <i>num_gid_entries</i> , <i>num_lid_entries</i> INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: <i>global_id</i> INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: <i>local_id</i> REAL(Zoltan_DOUBLE), INTENT(OUT), DIMENSION(*) :: <i>geom_vec</i> INTEGER(Zoltan_INT), INTENT(OUT) :: <i>ierr</i> < <i>type-data</i> > can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_ <i>x</i>) where <i>x</i> is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.

A **ZOLTAN_GEOM_FN** query function returns a vector of geometry values for a given object. The geometry vector is allocated by Zoltan to be of the size returned by a [ZOLTAN_NUM_GEOM_FN](#) query function.

Function Type:	ZOLTAN_GEOM_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>global_id</i>	The global ID of the object whose geometry values should be returned.
<i>local_id</i>	The local ID of the object whose geometry values should be returned.
<i>geom_vec</i>	Upon return, an array containing geometry values.
<i>ierr</i>	Error code to be set by function.

Graph-based Functions

C and C++:	typedef void ZOLTAN_NUM_EDGES_MULTI_FN (void * <i>data</i> , int <i>num_gid_entries</i> , int <i>num_lid_entries</i> , int <i>num_obj</i> , ZOLTAN_ID_PTR <i>global_ids</i> , ZOLTAN_ID_PTR <i>local_ids</i> , int * <i>num_edges</i> , int * <i>ierr</i>);
FORTRAN:	SUBROUTINE <i>Get_Num_Edges_Multi</i> (<i>data</i> , <i>num_gid_entries</i> , <i>num_lid_entries</i> , <i>num_obj</i> , <i>global_ids</i> , <i>local_ids</i> , <i>num_edges</i> , <i>ierr</i>) INTEGER(Zoltan_INT) :: <i>Get_Num_Edges</i>

```
<type-data>, INTENT(IN) :: data
INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries, num_obj
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_ids
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: local_ids
INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: num_edges
INTEGER(Zoltan_INT), INTENT(OUT) :: ierr
```

<type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_NUM_EDGES_MULTI_FN** query function returns the number of edges in the communication graph of the application for each object in a list of objects. That is, for each object in the *global_ids/local_ids* arrays, the number of objects with which the given object must share information is returned.

Function Type: **ZOLTAN_NUM_EDGES_MULTI_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>num_obj</i>	The number of object IDs in arrays <i>global_ids</i> and <i>local_ids</i> .
<i>global_ids</i>	Array of global IDs of objects whose number of edges should be returned.
<i>local_ids</i>	Array of local IDs of objects whose number of edges should be returned. (Optional.)
<i>num_edges</i>	Upon return, an array containing numbers of edges. For object <i>i</i> (specified by <i>global_ids[i*num_gid_entries]</i> and <i>local_ids[i*num_lid_entries]</i> , <i>i=0,1,...,num_obj-1</i>), the number of edges should be stored in <i>num_edges[i]</i> .
<i>ierr</i>	Error code to be set by function.

C and C++: typedef int **ZOLTAN_NUM_EDGES_FN** (void **data*, int *num_gid_entries*, int *num_lid_entries*, [ZOLTAN_ID_PTR](#) *global_id*, [ZOLTAN_ID_PTR](#) *local_id*, int **ierr*);

FORTTRAN: FUNCTION *Get_Num_Edges*(*data*, *num_gid_entries*, *num_lid_entries*, *global_id*, *local_id*, *ierr*)
INTEGER(Zoltan_INT) :: *Get_Num_Edges*
<type-data>, INTENT(IN) :: *data*
INTEGER(Zoltan_INT), INTENT(IN) :: *num_gid_entries*, *num_lid_entries*
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: *global_id*
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: *local_id*
INTEGER(Zoltan_INT), INTENT(OUT) :: *ierr*

<type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_NUM_EDGES_FN** query function returns the number of edges for a given object in the communication graph of the application (i.e., the number of objects with which the given object must share information).

Function Type: **ZOLTAN_NUM_EDGES_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>global_id</i>	The global ID of the object for which the number of edges should be returned.
<i>local_id</i>	The local ID of the object for which the number of edges should be returned.
<i>ierr</i>	Error code to be set by function.

Returned Value:

int	The number of edges for the object identified by <i>global_id</i> and <i>local_id</i> .
-----	-----------------------------------------------------------------------------------------

C and C++: typedef void **ZOLTAN_EDGE_LIST_MULTI_FN** (void **data*, int *num_gid_entries*,
int *num_lid_entries*, int *num_obj*, [ZOLTAN_ID_PTR](#) *global_ids*, [ZOLTAN_ID_PTR](#) *local_ids*,
int **num_edges*, [ZOLTAN_ID_PTR](#) *nbor_global_id*, int **nbor_procs*, int *wgt_dim*, float **ewgts*,
int **ierr*);

FORTRAN: SUBROUTINE *Get_Edge_List_Multi*(*data*, *num_gid_entries*, *num_lid_entries*, *num_obj*,
global_ids, *local_ids*, *num_edges*, *nbor_global_id*, *nbor_procs*, *wgt_dim*, *ewgts*, *ierr*)
 <*type-data*>, INTENT(IN) :: *data*
 INTEGER(Zoltan_INT), INTENT(IN) :: *num_gid_entries*, *num_lid_entries*, *num_obj*
 INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: *global_ids*
 INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: *local_ids*
 INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: *num_edges*
 INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: *nbor_global_id*
 INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: *nbor_procs*
 INTEGER(Zoltan_INT), INTENT(IN) :: *wgt_dim*
 REAL(Zoltan_FLOAT), INTENT(OUT), DIMENSION(*) :: *ewgts*
 INTEGER(Zoltan_INT), INTENT(OUT) :: *ierr*

 <*type-data*> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),
 DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x)
 where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_EDGE_LIST_MULTI_FN** query function returns lists of global IDs, processor IDs, and optionally edge weights for objects sharing edges with objects specified in the *global_ids* input array; objects share edges when they must share information with other objects. The arrays for the returned neighbor lists are allocated by Zoltan; their size is determined by a calls to [ZOLTAN_NUM_EDGES_MULTI_FN](#) or [ZOLTAN_NUM_EDGES_FN](#) query functions.

Function Type: **ZOLTAN_EDGE_LIST_MULTI_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>num_obj</i>	The number of object IDs in arrays <i>global_ids</i> and <i>local_ids</i> .

<i>global_ids</i>	Array of global IDs of objects whose edge lists should be returned.
<i>local_ids</i>	Array of local IDs of objects whose edge lists should be returned. (Optional.)
<i>num_edges</i>	An array containing numbers of edges for each object in <i>global_ids</i> . For object <i>i</i> (specified by <i>global_ids</i> [<i>i</i> * <i>num_gid_entries</i>] and <i>local_ids</i> [<i>i</i> * <i>num_lid_entries</i>], <i>i</i> =0,1,..., <i>num_obj</i> -1), the number of edges is stored in <i>num_edges</i> [<i>i</i>].
<i>nbor_global_id</i>	Upon return, an array of global IDs of objects sharing edges with the objects specified in <i>global_ids</i> . For object <i>i</i> (specified by <i>global_ids</i> [<i>i</i> * <i>num_gid_entries</i>] and <i>local_ids</i> [<i>i</i> * <i>num_lid_entries</i>], <i>i</i> =0,1,..., <i>num_obj</i> -1), edges are stored in <i>nbor_global_id</i> [<i>sum</i> * <i>num_gid_entries</i>] to <i>nbor_global_id</i> [(<i>sum</i> + <i>num_edges</i> [<i>i</i>])* <i>num_gid_entries</i> -1], where <i>sum</i> = the sum of <i>num_edges</i> [<i>j</i>] for <i>j</i> =0,1,..., <i>i</i> -1.
<i>nbor_procs</i>	Upon return, an array of processor IDs that identifies where the neighboring objects reside. For neighboring object <i>i</i> (stored in <i>nbor_global_id</i> [<i>i</i> * <i>num_gid_entries</i>]), the processor owning the neighbor is stored in <i>nbor_procs</i> [<i>i</i>].
<i>wgt_dim</i>	The number of weights associated with an edge (typically 1), or 0 if edge weights are not requested. This value is set through the parameter EDGE_WEIGHT_DIM .
<i>ewgts</i>	Upon return, an array of edge weights, where <i>ewgts</i> [<i>i</i> * <i>wgt_dim</i> :(<i>i</i> +1)* <i>wgt_dim</i> -1] corresponds to the weights for the <i>i</i> th edge. If <i>wgt_dim</i> =0, the return value of <i>ewgts</i> is undefined and may be NULL.
<i>ierr</i>	Error code to be set by function.

C and C++: `typedef void ZOLTAN_EDGE_LIST_FN (void *data, int num_gid_entries, int num_lid_entries, ZOLTAN_ID_PTR global_id, ZOLTAN_ID_PTR local_id, ZOLTAN_ID_PTR nbor_global_id, int *nbor_procs, int wgt_dim, float *ewgts, int *ierr);`

FORTTRAN: `SUBROUTINE Get_Edge_List(data, num_gid_entries, num_lid_entries, global_id, local_id, nbor_global_id, nbor_procs, wgt_dim, ewgts, ierr)`
`<type-data>, INTENT(IN) :: data`
`INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries`
`INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_id`
`INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: local_id`
`INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: nbor_global_id`
`INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: nbor_procs`
`INTEGER(Zoltan_INT), INTENT(IN) :: wgt_dim`
`REAL(Zoltan_FLOAT), INTENT(OUT), DIMENSION(*) :: ewgts`
`INTEGER(Zoltan_INT), INTENT(OUT) :: ierr`

`<type-data>` can be any of `INTEGER(Zoltan_INT)`, `DIMENSION(*)` or `REAL(Zoltan_FLOAT)`, `DIMENSION(*)` or `REAL(Zoltan_DOUBLE)`, `DIMENSION(*)` or `TYPE(Zoltan_User_Data_x)` where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_EDGE_LIST_FN** query function returns lists of global IDs, processor IDs, and optionally edge weights for objects sharing an edge with a given object (i.e., objects that must share information with the given object). The arrays for the returned neighbor lists are allocated by Zoltan; their size is determined by a call to [ZOLTAN_NUM_EDGES_MULTI_FN](#) or [ZOLTAN_NUM_EDGES_FN](#) query functions.

Function Type: **ZOLTAN_EDGE_LIST_FN_TYPE**

Arguments:

data Pointer to user-defined data.

<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>global_id</i>	The global ID of the object for which an edge list should be returned.
<i>local_id</i>	The local ID of the object for which an edge list should be returned.
<i>nbor_global_id</i>	Upon return, an array of global IDs of objects sharing edges with the given object.
<i>nbor_procs</i>	Upon return, an array of processor IDs that identifies where the neighboring objects reside.
<i>wgt_dim</i>	The number of weights associated with an edge (typically 1), or 0 if edge weights are not requested. This value is set through the parameter EDGE_WEIGHT_DIM .
<i>ewgts</i>	Upon return, an array of edge weights, where <i>ewgts</i> [<i>i</i> * <i>wgt_dim</i> :(<i>i</i> +1)* <i>wgt_dim</i> -1] corresponds to the weights for the <i>i</i> th edge. If <i>wgt_dim</i> =0, the return value of <i>ewgts</i> is undefined and may be NULL.
<i>ierr</i>	Error code to be set by function.

Hypergraph-based Functions

C and C++:	<pre>typedef void ZOLTAN_HG_SIZE_CS_FN (void *data, int *num_lists, int *num_pins, int *format, int *ierr);</pre>
FORTTRAN:	<pre>SUBROUTINE <i>Get_HG_Size_CS</i>(data, num_lists, num_pins, format, ierr) <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(OUT) :: num_lists INTEGER(Zoltan_INT), INTENT(OUT) :: num_pins INTEGER(Zoltan_INT), INTENT(OUT) :: format INTEGER(Zoltan_INT), INTENT(OUT) :: ierr</pre> <p><type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where <i>x</i> is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</p>

A hypergraph can be supplied to the Zoltan library in one of two compressed storage formats. Although hypergraphs are often used to represent the structure of sparse matrices, the Zoltan/PHG terminology is purely in terms of vertices and hyperedges (not rows or columns). The two compressed formats are analogous to CRS and CCS for matrices. In compressed hyperedge format (**ZOLTAN_COMPRESSED_EDGE**) a list of global hyperedge IDs is provided. Then a single list of the hypergraph pins, is provided. A pin is the connection between a vertex and a hyperedge (corresponds to a nonzero in a sparse matrix). Pins do not have separate IDs but are rather identified by the global ID of the vertex containing the pin, and implicitly also by the hyperedge ID. An example is provided below.

The other format is compressed vertex (**ZOLTAN_COMPRESSED_VERTEX**). In this format a list of vertex global IDs is provided. Then a list of pins ordered by vertex and then by hyperedge is provided. The pin ID in this case is the global ID of the hyperedge in which the pin appears. In both formats, an array must be provided pointing to the start in the list of pins where each hyperedge or vertex begins.

The purpose of this query function is to tell Zoltan in which format the application will supply the hypergraph, how

many vertices and hyperedges there will be, and how many pins. The actual hypergraph is supplied with a query function of the type [ZOLTAN_HG_CS_FN_TYPE](#).

This query function is required by all applications using the hypergraph methods of Zoltan (unless they are using the [graph-based](#) functions with hypergraph code instead).

Function Type:	ZOLTAN_HG_SIZE_CS_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>num_lists</i>	Upon return, the number of vertices (if using compressed vertex storage) or hyperedges (if using compressed hyperedge storage) that will be supplied to Zoltan by the application process.
<i>num_pins</i>	Upon return, the number of pins (connections between vertices and hyperedges) that will be supplied to Zoltan by the application process.
<i>format</i>	Upon return, the format in which the application process will provide the hypergraph to Zoltan. The options are ZOLTAN_COMPRESSED_EDGE and ZOLTAN_COMPRESSED_VERTEX .
<i>ierr</i>	Error code to be set by function.

C and C++:	<pre>typedef void ZOLTAN_HG_CS_FN (void *data, int num_gid_entries, int num_vtx_edge, int num_pins, int format, ZOLTAN_ID_PTR vtxedge_GID, int *vtxedge_ptr, ZOLTAN_ID_PTR pin_GID, int *ierr);</pre>
FORTRAN:	<pre>SUBROUTINE <i>Get_HG_CS</i>(data, num_gid_entries, num_vtx_edge, num_pins, format, vtxedge_GID, vtxedge_ptr, pin_GID, ierr) <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_vtx_edge, num_pins, format INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: vtxedge_GID INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: vtxedge_ptr INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: pin_GID INTEGER(Zoltan_INT), INTENT(OUT) :: ierr</pre> <p><type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where x is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</p>

A **ZOLTAN_HG_CS_FN** returns a hypergraph in a compressed storage (CS) format. The size and format of the data to be returned must have been supplied to Zoltan using a [ZOLTAN_HG_SIZE_CS_FN_TYPE](#) function.

When a hypergraph is distributed across multiple processes, Zoltan expects that all processes share a consistent global numbering scheme for hyperedges and vertices. Also, no two processes should return the same pin (matrix non-zero) in this query function. (Pin ownership is unique.)

This query function is required by all applications using the hypergraph methods of Zoltan (unless they are using the [graph-based](#) functions with hypergraph code instead).

Function Type:	ZOLTAN_HG_CS_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum

	value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_vtx_edge</i>	The number of global IDs that is expected to appear on return in <i>vtxedge_GID</i> . This may correspond to either vertices or (hyper-)edges.
<i>num_pins</i>	The number of pins that is expected to appear on return in <i>pin_GID</i> .
<i>format</i>	If <i>format</i> is ZOLTAN_COMPRESSED_EDGE , Zoltan expects that hyperedge global IDs will be returned in <i>vtxedge_GID</i> , and that vertex global IDs will be returned in <i>pin_GIDs</i> . If it is ZOLTAN_COMPRESSED_VERTEX , then vertex global IDs are expected to be returned in <i>vtxedge_GID</i> and hyperedge global IDs are expected to be returned in <i>pin_GIDs</i> .
<i>vtxedge_GID</i>	Upon return, a list of <i>num_vtx_edge</i> global IDs.
<i>vtxedge_ptr</i>	Upon return, this array contains <i>num_vtx_edge</i> integers such that the number of pins specified for hyperedge <i>j</i> (if <i>format</i> is ZOLTAN_COMPRESSED_EDGE) or vertex <i>j</i> (if <i>format</i> is ZOLTAN_COMPRESSED_VERTEX) is <i>vtxedge_ptr[j+1]-vtxedge_ptr[j]</i> . If <i>format</i> is ZOLTAN_COMPRESSED_EDGE , <i>vtxedge_ptr[j]*num_gid_entries</i> is the index into the array <i>pin_GID</i> where edge <i>j</i> 's pins (vertices belonging to edge <i>j</i>) begin; if <i>format</i> is ZOLTAN_COMPRESSED_VERTEX , <i>vtxedge_ptr[j]*num_gid_entries</i> is the index into the array <i>pin_GID</i> where vertex <i>j</i> 's pins (edges to which vertex <i>j</i> belongs) begin. Array indices begin at zero.
<i>pin_GID</i>	Upon return, a list of <i>num_pins</i> global IDs. This is the list of the pins contained in the hyperedges or vertices listed in <i>vtxedge_GID</i> .
<i>ierr</i>	Error code to be set by function.

Example

	vertex				
hyperedge	10	20	30	40	50
1	0	0	X	X	0
2	0	X	X	0	0
3	X	0	0	0	X

Compressed hyperedge storage:

$\text{vtxedge_GID} = \{1, 2, 3\}$
 $\text{vtxedge_ptr} = \{0, 2, 4\}$
 $\text{pin_GID} = \{30, 40, 20, 30, 10, 50\}$

Compressed vertex storage:

$\text{vtxedge_GID} = \{10, 20, 30, 40, 50\}$
 $\text{vtxedge_ptr} = \{0, 1, 2, 4, 5\}$
 $\text{pin_GID} = \{3, 2, 1, 2, 1, 3\}$

C and C++: typedef void **ZOLTAN_HG_SIZE_EDGE_WTS_FN** (void *data, int *num_edges, int *ierr);

FORTRAN: SUBROUTINE *Get_HG_Size_Edge_Wts*(data, num_edges, ierr)

<type-data>, INTENT(IN) :: data

INTEGER(Zoltan_INT), INTENT(OUT) :: num_edges

INTEGER(Zoltan_INT), INTENT(OUT) :: ierr

<type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),

DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x)

where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_HG_SIZE_EDGE_WTS_FN** returns the number of hyperedges for which a process will supply edge weights. The number of weights per hyperedge was supplied by the application with the [EDGE_WEIGHT_DIM](#) parameter. The actual edge weights will be supplied with a [ZOLTAN_HG_EDGE_WTS_FN_TYPE](#) function.

This query function is not required. If no hyperedge weights are supplied, Zoltan will assume every hyperedge has weight 1.0.

Function Type: **ZOLTAN_HG_SIZE_EDGE_WTS_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_edges</i>	Upon return, the number of hyperedges for which edge weights will be supplied.
<i>ierr</i>	Error code to be set by function.

C and C++:	<pre>typedef void ZOLTAN_HG_EDGE_WTS_FN (void *data, int num_gid_entries, int num_lid_entries, int num_edges, int edge_weight_dim, ZOLTAN_ID_PTR edge_GID, ZOLTAN_ID_PTR edge_LID, float *edge_weight, int *ierr);</pre>
FORTRAN:	<pre>SUBROUTINE <i>Get_HG_Edge_Wts</i>(data, num_gid_entries, num_lid_entries, num_edges, edge_weight_dim, edge_GID, edge_LID, edge_weight, ierr) <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries, num_edges, edge_weight_dim INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: edge_GID INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: edge_LID REAL(Zoltan_FLOAT), INTENT(OUT), DIMENSION(*) :: edge_weight INTEGER(Zoltan_INT), INTENT(OUT) :: ierr</pre> <p><type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where x is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</p>

A **ZOLTAN_HG_EDGE_WTS_FN** returns edges weights for a set of hypergraph edges. The number of weights supplied for each hyperedge should equal the value of the [EDGE_WEIGHT_DIM](#) parameter. In the case of a hypergraph which is distributed across multiple processes, if more than one process supplies edge weights for the same hyperedge, the different edge weights will be resolved according to the value of the [PHG_EDGE_WEIGHT_OPERATION](#) parameter.

This query function is not required. If no hyperedge weights are supplied, Zoltan will assume every hyperedge has weight 1.0.

Function Type: **ZOLTAN_HG_SIZE_EDGE_WTS_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>num_edges</i>	The number of hyperedges for which edge weights should be supplied in the <i>edge_weight</i> array.
<i>edge_weight_dim</i>	The number of weights which should be supplied for each hyperedge. This is also the value of the EDGE_WEIGHT_DIM parameter.
<i>edge_GID</i>	Upon return, this array should contain the global IDs of the <i>num_edges</i> hyperedges for which

the application is supplying edge weights.

<i>edge_LID</i>	Upon return, this array can optionally contain the local IDs of the <i>num_edges</i> hyperedges for which the application is supplying edge weights.
<i>edge_weight</i>	Upon return, this array should contain the weights for each edge listed in the <i>edge_GID</i> . If <i>edge_weight_dim</i> is greater than one, all weights for one hyperedge are listed before the weights for the next hyperedge are listed.
<i>ierr</i>	Error code to be set by function.

C and C++:	typedef int ZOLTAN_NUM_FIXED_OBJ_FN (void * <i>data</i> , int * <i>ierr</i>);
FORTTRAN:	FUNCTION <i>Get_Num_Fixed_Obj</i> (<i>data</i> , <i>ierr</i>) INTEGER(Zoltan_INT) :: Get_Num_Fixed_Obj < <i>type-data</i> >, INTENT(IN) :: <i>data</i> INTEGER(Zoltan_INT), INTENT(OUT) :: <i>ierr</i> < <i>type-data</i> > can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_ <i>x</i>) where <i>x</i> is 1, 2, 3 or 4. See the section on Fortran

A **ZOLTAN_NUM_FIXED_OBJ_FN** returns the number of objects on a given processor fixed to particular parts. These objects will not be assigned to a part other than the user specified part by the parallel hypergraph algorithm.

This query function is not required. If it is not defined, all objects are candidates for migration to new parts.

Function Type:	ZOLTAN_NUM_FIXED_OBJ_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>ierr</i>	Error code to be set by function.
Returned Value:	
int	The number of objects on this processor that are to be fixed to a specific part.

C and C++:	typedef void ZOLTAN_FIXED_OBJ_LIST_FN (void * <i>data</i> , int <i>num_fixed_obj</i> , int <i>num_gid_entries</i> , ZOLTAN_ID_PTR <i>fixed_gids</i> , int * <i>fixed_parts</i> , int * <i>ierr</i>);
FORTTRAN:	SUBROUTINE <i>Get_Fixed_Obj_List</i> (<i>data</i> , <i>num_fixed_obj</i> , <i>num_gid_entries</i> , <i>fixed_gids</i> , <i>fixed_parts</i> , <i>ierr</i>) < <i>type-data</i> >, INTENT(IN) :: <i>data</i> INTEGER(Zoltan_INT), INTENT(IN) :: <i>num_fixed_obj</i> INTEGER(Zoltan_INT), INTENT(IN) :: <i>num_gid_entries</i> INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: <i>fixed_gids</i> INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: <i>fixed_parts</i> INTEGER(Zoltan_INT), INTENT(OUT) :: <i>ierr</i> < <i>type-data</i> > can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_ <i>x</i>) where <i>x</i> is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.

A **ZOLTAN_FIXED_OBJ_LIST_FN** query function fills two arrays with information about the objects that should not be moved from their user assigned parts. These arrays are allocated (and subsequently freed) by Zoltan; their size is determined by a call to a [ZOLTAN_NUM_FIXED_OBJ_FN](#) query function to get the array size.

A process should only fix the part of objects that it *owns*, that is, objects that it had previously supplied in a [ZOLTAN_OBJ_LIST_FN](#) query function. It is an error to list the global ID of an object owned by another process.

Function Type:	ZOLTAN_FIXED_OBJ_LIST_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>num_fixed_obj</i>	The number of objects you will list in the two output arrays.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>fixed_gids</i>	Upon return, an array of unique global IDs for all objects assigned to the processor which are to be fixed to a part.
<i>fixed_parts</i>	Upon return, an array of part numbers, one for each object listed in the global ID array. These objects will not be migrated from this assigned part.
<i>ierr</i>	Error code to be set by function.

Tree-based Functions

C and C++:	typedef int ZOLTAN_NUM_COARSE_OBJ_FN (void * <i>data</i> , int * <i>ierr</i>);
FORTRAN:	<pre>FUNCTION <i>Get_Num_Coarse_Obj</i>(<i>data</i>, <i>ierr</i>) INTEGER(Zoltan_INT) :: Get_Num_Coarse_Obj <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(OUT) :: ierr</pre> <p><type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where <i>x</i> is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</p>

A **ZOLTAN_NUM_COARSE_OBJ_FN** query function returns the number of objects (elements) in the initial coarse grid.

Function Type:	ZOLTAN_NUM_COARSE_OBJ_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>ierr</i>	Error code to be set by function.
Returned Value:	
int	The number of objects in the coarse grid.

C and C++:	typedef void ZOLTAN_COARSE_OBJ_LIST_FN (void * <i>data</i> , int <i>num_gid_entries</i> , int <i>num_lid_entries</i> , ZOLTAN_ID_PTR <i>global_ids</i> , ZOLTAN_ID_PTR <i>local_ids</i> , int * <i>assigned</i> ,
------------	----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------


```
int *num_vert, ZOLTAN\_ID\_PTR vertices, int *in_order, ZOLTAN\_ID\_PTR in_vertex,
ZOLTAN\_ID\_PTR out_vertex, int *ierr);
```

FORTTRAN: SUBROUTINE *Get_Coarse_Obj_List*(data, num_gid_entries, num_lid_entries, global_ids, local_ids, assigned, num_vert, vertices, in_order, in_vertex, out_vertex, ierr)
 <type-data>, INTENT(IN) :: data
 INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries
 INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: global_ids
 INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: local_ids
 INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: assigned, num_vert, vertices,
 in_vertex, out_vertex
 INTEGER(Zoltan_INT), INTENT(OUT) :: in_order, ierr

<type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_COARSE_OBJ_LIST_FN** query function returns lists of global IDs, local IDs, vertices, and order information for all objects (elements) of the initial coarse grid. The vertices are designated by a global ID such that if two elements share a vertex then the same ID designates that vertex in both elements and on all processors. The user may choose to provide the order in which the elements should be traversed or have Zoltan determine the order. If the user provides the order, then entry and exit vertices for a path through the elements may also be provided. The arrays for the returned values are allocated by Zoltan; their size is determined by a call to a [ZOLTAN_NUM_COARSE_OBJ_FN](#) query function.

Function Type: **ZOLTAN_COARSE_OBJ_LIST_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>global_ids</i>	Upon return, an array of global IDs of all objects in the coarse grid.
<i>local_ids</i>	Upon return, an array of local IDs of all objects in the coarse grid. (Optional.)
<i>assigned</i>	Upon return, an array of integers indicating whether or not each object is currently assigned to this processor. A value of 1 indicates it is assigned to this processor; a value of 0 indicates it is assigned to some other processor. For elements that have been refined, it is ignored unless weights are assigned to interior nodes of the tree.
<i>num_vert</i>	Upon return, an array containing the number of vertices for each object.
<i>vertices</i>	Upon return, an array of global IDs of the vertices of each object. If the number of vertices for objects 0 through <i>i</i> -1 is <i>N</i> , then the vertices for object <i>i</i> are in <i>vertices</i> [<i>N</i> * <i>num_gid_entries</i> : (<i>N</i> + <i>num_vert</i> [<i>i</i>])* <i>num_gid_entries</i>]
<i>in_order</i>	Upon return, 1 if the user is providing the objects in the order in which they should be traversed, or 0 if Zoltan should determine the order.
<i>in_vertex</i>	Upon return, an array of global IDs of the vertices through which to enter each element in the user provided traversal. It is required only if the user is providing the order for the coarse grid objects (i.e., <i>in_order</i> ==1) and allowing Zoltan to select the order of the children in at least one invocation of ZOLTAN_CHILD_LIST_FN .
<i>out_vertex</i>	Upon return, an array of global IDs of the vertex through which to exit each element in the user provided traversal. The same provisions hold as for <i>in_vertex</i> .

<i>ierr</i>	Error code to be set by function.
<hr/>	
C and C++:	<pre>typedef int ZOLTAN_FIRST_COARSE_OBJ_FN (void *data, int num_gid_entries, int num_lid_entries, ZOLTAN_ID_PTR global_id, ZOLTAN_ID_PTR local_id, int *assigned, int *num_vert, ZOLTAN_ID_PTR vertices, int *in_order, ZOLTAN_ID_PTR in_vertex, ZOLTAN_ID_PTR out_vertex, int *ierr);</pre>
FORTTRAN:	<pre>FUNCTION <i>Get_First_Coarse_Obj</i>(data, num_gid_entries, num_lid_entries, global_id, local_id, assigned, num_vert, vertices, in_order, in_vertex, out_vertex, ierr) INTEGER(Zoltan_INT) :: Get_First_Coarse_Obj <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: global_id INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: local_id INTEGER(Zoltan_INT), INTENT(OUT) :: assigned, num_vert, in_order, ierr INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: vertices, in_vertex, out_vertex <type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where x is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</pre>

A **ZOLTAN_FIRST_COARSE_OBJ_FN** query function initializes an iteration over the objects of the initial coarse grid. It returns the global ID, local ID, vertices, and order information for the first object (element) of the initial coarse grid. Subsequent calls to a [ZOLTAN_NEXT_COARSE_OBJ_FN](#) iterate over and return other objects from the coarse grid. The vertices are designated by a global ID such that if two elements share a vertex then the same ID designates that vertex in both elements and on all processors. The user may choose to provide the order in which the elements should be traversed, or have Zoltan determine the order. If the user provides the order, then entry and exit vertices for a path through the elements may also be provided.

Function Type:	ZOLTAN_FIRST_COARSE_OBJ_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>global_ids</i>	Upon return, the global ID of the first object in the coarse grid.
<i>local_ids</i>	Upon return, the local ID of the first object in the coarse grid. (Optional.)
<i>assigned</i>	Upon return, an integer indicating whether or not this object is currently assigned to this processor. A value of 1 indicates it is assigned to this processor; a value of 0 indicates it is assigned to some other processor. For elements that have been refined, it is ignored unless weights are assigned to interior nodes of the tree.
<i>num_vert</i>	Upon return, the number of vertices for this object.
<i>vertices</i>	Upon return, an array of global IDs of the vertices of this object.
<i>in_order</i>	Upon return, 1 if the user is providing the objects in the order in which they should be traversed, or 0 if Zoltan should determine the order.
<i>in_vertex</i>	Upon return, the vertex through which to enter this element in the user provided traversal. It is required only if the user is providing the order for the coarse grid objects (i.e., <i>in_order</i> ==1) and allowing Zoltan to select the order of the children in at least one invocation of

[ZOLTAN_CHILD_LIST_FN.](#)

<i>out_vertex</i>	Upon return, the vertex through which to exit this element in the user provided traversal. The same provisions hold as for <i>in_vertex</i> .
<i>ierr</i>	Error code to be set by function.
Returned Value:	
1	If <i>global_id</i> and <i>local_id</i> contain valid IDs of the first object in the coarse grid.
0	If no coarse grid is available.

C and C++:	<pre>typedef int ZOLTAN_NEXT_COARSE_OBJ_FN (void *data, int num_gid_entries, int num_lid_entries, ZOLTAN_ID_PTR global_id, ZOLTAN_ID_PTR local_id, ZOLTAN_ID_PTR next_global_id, ZOLTAN_ID_PTR next_local_id, int *assigned, int *num_vert, ZOLTAN_ID_PTR vertices, ZOLTAN_ID_PTR in_vertex, ZOLTAN_ID_PTR out_vertex, int *ierr);</pre>
FORTTRAN:	<pre>FUNCTION <i>Get_Next_Coarse_Obj</i>(data, num_gid_entries, num_lid_entries, global_id, local_id, next_global_id, next_local_id, assigned, num_vert, vertices, in_vertex, out_vertex, ierr) INTEGER(Zoltan_INT) :: Get_Next_Coarse_Obj <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_id INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: local_id INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: next_global_id INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: next_local_id INTEGER(Zoltan_INT), INTENT(OUT) :: assigned, num_vertex, ierr INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: vertices, in_vertex, out_vertex <type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where x is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</pre>

A **ZOLTAN_NEXT_COARSE_OBJ_FN** query function is an iterator function that returns the next object in the initial coarse grid. The first object of the iteration is provided by a [ZOLTAN_FIRST_COARSE_OBJ_FN](#) query function.

Function Type: **ZOLTAN_NEXT_COARSE_OBJ_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>global_id</i>	The global ID of the previous object in the coarse grid.
<i>local_id</i>	The local ID of the previous object in the coarse grid.
<i>next_global_id</i>	Upon return, the global ID of the next object in the coarse grid.
<i>next_local_id</i>	Upon return, the local ID of the next object in the coarse grid.
<i>assigned</i>	Upon return, an integer indicating whether or not this object is currently assigned to this processor. A value of 1 indicates it is assigned to this processor; a value of 0 indicates it is assigned to some other processor. For elements that have been refined, it is ignored unless

weights are assigned to interior nodes of the tree.

<i>num_vert</i>	Upon return, the number of vertices for this object.
<i>vertices</i>	Upon return, an array of global IDs of the vertices of this object.
<i>in_vertex</i>	Upon return, the vertex through which to enter this element in the user provided traversal. It is required only if the user is providing the order for the coarse grid objects (i.e., <i>in_order</i> ==1) and allowing Zoltan to select the order of the children in at least one invocation of ZOLTAN_CHILD_LIST_FN .
<i>out_vertex</i>	Upon return, the vertex through which to exit this element in the user provided traversal. The same provisions hold as for <i>in_vertex</i> .
<i>ierr</i>	Error code to be set by function.

Returned Value:

1	If <i>global_id</i> and <i>local_id</i> contain valid IDs of the next object in the coarse grid.
0	If no more objects are available.

C and C++:	typedef int ZOLTAN_NUM_CHILD_FN (void * <i>data</i> , int <i>num_gid_entries</i> , int <i>num_lid_entries</i> , ZOLTAN_ID_PTR <i>global_id</i> , ZOLTAN_ID_PTR <i>local_id</i> , int * <i>ierr</i>);
FORTTRAN:	FUNCTION <i>Get_Num_Child</i> (<i>data</i> , <i>num_gid_entries</i> , <i>num_lid_entries</i> , <i>global_id</i> , <i>local_id</i> , <i>ierr</i>) INTEGER(Zoltan_INT) :: Get_Num_Child < <i>type-data</i> >, INTENT(IN) :: <i>data</i> INTEGER(Zoltan_INT), INTENT(IN) :: <i>num_gid_entries</i> , <i>num_lid_entries</i> INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: <i>global_id</i> INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: <i>local_id</i> INTEGER(Zoltan_INT), INTENT(OUT) :: <i>ierr</i> < <i>type-data</i> > can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_ <i>x</i>) where <i>x</i> is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.

A **ZOLTAN_NUM_CHILD_FN** query function returns the number of children of the element with the given global and local IDs. If the element has not been refined, the number of children is 0.

Function Type: **ZOLTAN_NUM_CHILD_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>global_id</i>	The global ID of the object for which the number of children is requested.
<i>local_id</i>	The local ID of the object for which the number of children is requested.
<i>ierr</i>	Error code to be set by function.

Returned Value:

int	The number of children.
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C and C++:	typedef void ZOLTAN_CHILD_LIST_FN (void * <i>data</i> , int <i>num_gid_entries</i> , int <i>num_lid_entries</i> ,
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[ZOLTAN_ID_PTR](#) *parent_gid*, [ZOLTAN_ID_PTR](#) *parent_lid*, [ZOLTAN_ID_PTR](#) *child_gids*,
[ZOLTAN_ID_PTR](#) *child_lids*, int **assigned*, int **num_vert*, [ZOLTAN_ID_PTR](#) *vertices*,
[ZOLTAN_REF_TYPE](#) **ref_type*, [ZOLTAN_ID_PTR](#) *in_vertex*, [ZOLTAN_ID_PTR](#) *out_vertex*,
int **ierr*);

FORTTRAN: SUBROUTINE *Get_Child_List*(*data*, *num_gid_entries*, *num_lid_entries*, *parent_gid*, *parent_lid*,
child_gids, *child_lids*, *assigned*, *num_vert*, *vertices*, *ref_type*, *in_vertex*, *out_vertex*, *ierr*)
<*type-data*>, INTENT(IN) :: *data*
INTEGER(Zoltan_INT), INTENT(IN) :: *num_gid_entries*, *num_lid_entries*
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: *parent_gid*
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: *parent_lid*
INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: *child_gids*
INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: *child_lids*
INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: *assigned*, *num_vert*, *vertices*,
in_vertex, *out_vertex*
INTEGER(Zoltan_INT), INTENT(OUT) :: *ref_type*, *ierr*

<*type-data*> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),
DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x)
where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_CHILD_LIST_FN** query function returns lists of global IDs, local IDs, vertices, and order information for all children of a refined element. The vertices are designated by a global ID such that if two elements share a vertex then the same ID designates that vertex in both elements and on all processors. The user may choose to provide the order in which the children should be traversed, or have Zoltan determine the order based on the type of element refinement used to create the children. If the user provides the order, then entry and exit vertices for a path through the elements may also be provided. The arrays for the returned values are allocated by Zoltan; their size is determined by a call to a [ZOLTAN_NUM_CHILD_FN](#) query function.

Function Type: ZOLTAN_CHILD_LIST_FN_TYPE

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>parent_gid</i>	The global ID of the object whose children are requested.
<i>parent_lid</i>	The local ID of the object whose children are requested.
<i>child_gids</i>	Upon return, an array of global IDs of all children of this object.
<i>child_lids</i>	Upon return, an array of local IDs of all children of this object. (Optional.)
<i>assigned</i>	Upon return, an array of integers indicating whether or not each child is currently assigned to this processor. A value of 1 indicates it is assigned to this processor; a value of 0 indicates it is assigned to some other processor. For children that have been further refined, it is ignored unless weights are assigned to interior nodes of the tree.
<i>num_vert</i>	Upon return, an array containing the number of vertices for each object.
<i>vertices</i>	Upon return, an array of global IDs of the vertices of each object. If the number of vertices for objects 0 through <i>i</i> -1 is <i>N</i> , then the vertices for object <i>i</i> are in <i>vertices</i> [<i>N*num_gid_entries</i> : (<i>N+num_vert</i> [<i>i</i>])* <i>num_gid_entries</i>]
<i>ref_type</i>	Upon return, a value indicating what type of refinement was used to create the children. This determines how the children will be ordered. The values currently supported are:

ZOLTAN_TRI_BISECT Bisection of triangles.
ZOLTAN_QUAD_QUAD Quadrisection of quadrilaterals.
ZOLTAN_HEX3D_OCT Octasection of hexahedra.
ZOLTAN_OTHER_REF All other forms of refinement.
ZOLTAN_IN_ORDER Traverse the children in the order in which they are provided.

<i>in_vertex</i>	Upon return, an array of global IDs of the vertex through which to enter each element in the user provided traversal. It is required only if the user is providing the order for the children of this element (i.e., <i>ref_type</i> == <i>ZOLTAN_IN_ORDER</i>) but does not provide the order for the children of at least one of those children.
<i>out_vertex</i>	Upon return, an array of global IDs of the vertex through which to exit each element in the user provided traversal. The same provisions hold as for <i>in_vertex</i> .
<i>ierr</i>	Error code to be set by function.

C and C++:	<pre>typedef void ZOLTAN_CHILD_WEIGHT_FN (void *data, int num_gid_entries, int num_lid_entries, ZOLTAN_ID_PTR global_id, ZOLTAN_ID_PTR local_id, int wgt_dim, float *obj_wgt, int *ierr);</pre>
FORTTRAN:	<pre>SUBROUTINE <i>Get_Child_Weight</i>(data, num_gid_entries, num_lid_entries, global_id, local_id, wgt_dim, obj_wgt, ierr) <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_id INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: local_id INTEGER(Zoltan_INT), INTENT(IN) :: wgt_dim REAL(Zoltan_FLOAT), INTENT(OUT), DIMENSION(*) :: obj_wgt INTEGER(Zoltan_INT), INTENT(OUT) :: ierr</pre> <p><type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where x is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</p>

A **ZOLTAN_CHILD_WEIGHT_FN** query function returns the weight of an object. Interior nodes of the refinement tree as well as the leaves are allowed to have weights.

Function Type: **ZOLTAN_CHILD_WEIGHT_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES . (It should be zero if local ids are not used.)
<i>global_id</i>	The global ID of the object whose weight is requested.
<i>local_id</i>	The local ID of the object whose weight is requested.
<i>wgt_dim</i>	The number of weights associated with an object (typically 1), or 0 if weights are not requested. This value is set through the parameter OBJ_WEIGHT_DIM .
<i>obj_wgt</i>	Upon return, an array containing the object's weights. If <i>wgt_dim</i> =0, the return value of <i>obj_wgts</i> is undefined and may be NULL.
<i>ierr</i>	Error code to be set by function.

Hierarchical Partitioning Functions (used only by method HIER)

C and C++: typedef int **ZOLTAN_HIER_NUM_LEVELS_FN** (void **data*, int **ierr*);
FORTRAN: FUNCTION *Get_Hier_Num_Levels*(*data*, *nbor_proc*, *ierr*)
 INTEGER(Zoltan_INT) :: Get_Hier_Num_Levels
 <*type-data*>, INTENT(IN) :: *data*
 INTEGER(Zoltan_INT), INTENT(OUT) :: *ierr*

 <*type-data*> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),
 DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x)
 where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_HIER_NUM_LEVELS_FN** query function returns, for the calling processor, the number of levels of hierarchy for hierarchical load balancing.

Function Type:	ZOLTAN_HIER_NUM_LEVELS_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>ierr</i>	Error code to be set by function.
Returned Value:	
int	the number of levels of balancing hierarchy for method HIER.

C and C++: typedef int **ZOLTAN_HIER_PART_FN** (void **data*, int *level*, int **ierr*);
FORTRAN: FUNCTION *Get_Hier_Part*(*data*, *level*, *ierr*)
 INTEGER(Zoltan_INT) :: Get_Hier_Part
 <*type-data*>, INTENT(IN) :: *data*
 INTEGER(Zoltan_INT), INTENT(IN) :: *level*
 INTEGER(Zoltan_INT), INTENT(OUT) :: *ierr*

 <*type-data*> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),
 DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x)
 where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_HIER_PART_FN** query function gets the part number to be used for the given level of a hierarchical balancing procedure.

Function Type:	ZOLTAN_HIER_PART_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>level</i>	The level of a hierarchical balancing for which the part ID is requested.
<i>ierr</i>	Error code to be set by function.
Returned Value:	
int	The part number the process is to compute for this level.

C and C++:	typedef void ZOLTAN_HIER_METHOD_FN (void * <i>data</i> , int <i>level</i> , struct Zoltan_Struct * <i>zz</i> , int * <i>ierr</i>);
FORTTRAN:	SUBROUTINE <i>Get_Hier_Method</i> (<i>data</i> , <i>level</i> , <i>zz</i> , <i>ierr</i>) < <i>type-data</i> >, INTENT(IN) :: <i>data</i> INTEGER(Zoltan_INT), INTENT(IN) :: <i>level</i> TYPE(Zoltan_Struct), INTENT(IN) :: <i>zz</i> INTEGER(Zoltan_INT), INTENT(OUT) :: <i>ierr</i>

<*type-data*> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_*x*) where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_HIER_METHOD_FN** query function provides to the calling process the Zoltan_Struct to be used to guide the partitioning and load balancing at the given level in the hierarchy. This Zoltan_Struct can be passed to Zoltan_Set_Param to set load balancing parameters for this level in the hierarchical balancing.

Function Type:	ZOLTAN_HIER_METHOD_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>level</i>	Level in the hierarchy being considered.
<i>zz</i>	Zoltan_Struct to use to set parameters.
<i>ierr</i>	Error code to be set by function.

Migration Query Functions

The following query functions must be registered to use any of the migration tools described in [Migration Functions](#). These functions should NOT use interprocessor communication.

[ZOLTAN_OBJ_SIZE_MULTI_FN](#) or [ZOLTAN_OBJ_SIZE_FN](#)
[ZOLTAN_PACK_OBJ_MULTI_FN](#) or [ZOLTAN_PACK_OBJ_FN](#)
[ZOLTAN_UNPACK_OBJ_MULTI_FN](#) or [ZOLTAN_UNPACK_OBJ_FN](#)

The "MULTI_" versions of the packing/unpacking functions take lists of IDs as input and pack/unpack data for all objects in the lists. Only one function of each type must be provided (e.g., either a [ZOLTAN_PACK_OBJ_FN](#) or [ZOLTAN_PACK_OBJ_MULTI_FN](#), but not both).

Optional, additional query functions for migration may also be registered; these functions are called at the beginning, middle, and end of migration in [Zoltan_Migrate](#). Interprocessor communication is allowed in these functions.

[ZOLTAN_PRE_MIGRATE_PP_FN](#)
[ZOLTAN_MID_MIGRATE_PP_FN](#)
[ZOLTAN_POST_MIGRATE_PP_FN](#)

For [backward compatibility](#) with previous versions of Zoltan, the following functions may be used with [Zoltan_Help_Migrate](#).

[ZOLTAN_PRE_MIGRATE_FN](#)
[ZOLTAN_MID_MIGRATE_FN](#)
[ZOLTAN_POST_MIGRATE_FN](#)

C and C++:	<pre>typedef int ZOLTAN_OBJ_SIZE_FN(void *data, int num_gid_entries, int num_lid_entries, ZOLTAN_ID_PTR global_id, ZOLTAN_ID_PTR local_id, int *ierr);</pre>
FORTTRAN:	<pre>FUNCTION <i>Obj_Size</i>(data, num_gid_entries, num_lid_entries, global_id, local_id, ierr) INTEGER(Zoltan_INT) :: Obj_Size <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_id, local_id INTEGER(Zoltan_INT), INTENT(OUT) :: ierr <type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where x is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</pre>

A **ZOLTAN_OBJ_SIZE_FN** query function returns the size (in bytes) of the data buffer that is needed to pack all of a single object's data.

Function Type: **ZOLTAN_OBJ_SIZE_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES .
<i>global_id</i>	Pointer to the global ID of the object.
<i>local_id</i>	Pointer to the local ID of the object.
<i>ierr</i>	Error code to be set by function.

Returned Value:

int	The size (in bytes) of the required data buffer.
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C and C++:	<pre>typedef void ZOLTAN_OBJ_SIZE_MULTI_FN (void *data, int num_gid_entries, int num_lid_entries, int num_ids, ZOLTAN_ID_PTR global_ids, ZOLTAN_ID_PTR local_ids, int *sizes, int *ierr);</pre>
FORTTRAN:	<pre>SUBROUTINE <i>Obj_Size_Multi</i>(data, num_gid_entries, num_lid_entries, num_ids, global_ids, local_ids, sizes, ierr) <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries, num_ids INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_ids, local_ids INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: sizes INTEGER(Zoltan_INT), INTENT(OUT) :: ierr</pre> <p><type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where <i>x</i> is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</p>

A **ZOLTAN_OBJ_SIZE_MULTI_FN** query function is the multiple-ID version of [ZOLTAN_OBJ_SIZE_FN](#). For a list of objects, it returns the per-objects sizes (in bytes) of the data buffers needed to pack object data.

Function Type: **ZOLTAN_OBJ_SIZE_MULTI_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES .
<i>num_ids</i>	The number of objects whose sizes are to be returned.
<i>global_ids</i>	An array of global IDs of the objects. The ID for the <i>i</i> -th object begins in <i>global_ids[i*num_gid_entries]</i> .
<i>local_ids</i>	An array of local IDs of the objects. The ID for the <i>i</i> -th object begins in <i>local_ids[i*num_lid_entries]</i> .

sizes Upon return, array of sizes (in bytes) for each object in the ID lists.
ierr Error code to be set by function.

Returned Value:

int The size (in bytes) of the required data buffer.

C and C++:	<pre>typedef void ZOLTAN_PACK_OBJ_FN (void *data, int num_gid_entries, int num_lid_entries, ZOLTAN_ID_PTR global_id, ZOLTAN_ID_PTR local_id, int dest, int size, char *buf, int *ierr);</pre>
FORTTRAN:	<pre>SUBROUTINE <i>Pack_Obj</i>(data, num_gid_entries, num_lid_entries, global_id, local_id, dest, size, buf, ierr) <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_id INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: local_id INTEGER(Zoltan_INT), INTENT(IN) :: dest, size INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: buf INTEGER(Zoltan_INT), INTENT(OUT) :: ierr</pre> <p><type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where <i>x</i> is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</p>

A **ZOLTAN_PACK_OBJ_FN** query function allows the application to tell Zoltan how to copy all needed data for a given object into a communication buffer. The object's data can then be sent to another processor as part of data migration. It may also perform other operations, such as removing the object from the processor's data structure. This routine is called by [Zoltan_Migrate](#) for each object to be sent to another processor.

Function Type: **ZOLTAN_PACK_OBJ_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES .
<i>global_id</i>	The global ID of the object for which data should be copied into the communication buffer.
<i>local_id</i>	The local ID of the object for which data should be copied into the communication buffer.
<i>dest</i>	The destination part (i.e., the part to which the object is being sent)
<i>size</i>	The size (in bytes) of the communication buffer for the specified object. This value is at least as large as the value returned by the ZOLTAN_OBJ_SIZE_MULT_FN or ZOLTAN_OBJ_SIZE_FN query function; it may be slightly larger due to padding for data alignment in the buffer.
<i>buf</i>	The starting address of the communication buffer into which the object's data should be

packed.
ierr Error code to be set by function.

C and C++:	<pre>typedef void ZOLTAN_PACK_OBJ_MULTI_FN (void *data, int num_gid_entries, int num_lid_entries, int num_ids, ZOLTAN_ID_PTR global_ids, ZOLTAN_ID_PTR local_ids, int *dest, int *sizes, int *idx, char *buf, int *ierr);</pre>
FORTTRAN:	<pre>SUBROUTINE <i>Pack_Obj_Multi</i>(data, num_gid_entries, num_lid_entries, num_ids, global_ids, local_ids, dest, sizes, idx, buf, ierr) <type-data>, INTENT(IN) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries, num_ids INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_ids INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: local_ids INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: dest INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: sizes INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: idx INTEGER(Zoltan_INT), INTENT(OUT), DIMENSION(*) :: buf INTEGER(Zoltan_INT), INTENT(OUT) :: ierr</pre> <p><type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where <i>x</i> is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</p>

A **ZOLTAN_PACK_OBJ_MULTI_FN** query function is the multiple-ID version of a [ZOLTAN_PACK_OBJ_FN](#). It allows the application to tell Zoltan how to copy all needed data for a given list of objects into a communication buffer.

Function Type:	ZOLTAN_PACK_OBJ_FN_MULTI_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES .
<i>num_ids</i>	The number of objects to be packed.
<i>global_ids</i>	An array of global IDs of the objects. The ID for the <i>i</i> -th object begins in <i>global_ids[i*num_gid_entries]</i> .
<i>local_ids</i>	An array of local IDs of the objects. The ID for the <i>i</i> -th object begins in <i>local_ids[i*num_lid_entries]</i> .
<i>dest</i>	An array of destination part numbers (i.e., the parts to which the objects are being sent)
<i>sizes</i>	An array containing the per-object sizes (in bytes) of the communication buffer for each

object. Each value is at least as large as the corresponding value returned by the [ZOLTAN_OBJ_SIZE_MULTI_FN](#) or [ZOLTAN_OBJ_SIZE_FN](#) query function; it may be slightly larger due to padding for data alignment in the buffer.

<i>idx</i>	For each object, an index into the <i>buf</i> array giving the starting location of that object's data. Data for the <i>i</i> -th object are stored in <i>buf[idx[i]]</i> , <i>buf[idx[i]+1]</i> , ..., <i>buf[idx[i]+sizes[i]-1]</i> . Because Zoltan adds some tag information to packed data, <i>idx[i] != sum[j=0,i-1](sizes[j])</i> .
<i>buf</i>	The address of the communication buffer into which the objects' data should be packed.
<i>ierr</i>	Error code to be set by function.

```
C and C++:      typedef void ZOLTAN_UNPACK_OBJ_FN (  
                void *data,  
                int num_gid_entries,  
                ZOLTAN\_ID\_PTR global_id,  
                int size,  
                char *buf,  
                int *ierr);
```

```
FORTRAN:      SUBROUTINE Unpack_Obj(data, num_gid_entries, global_id, size, buf, ierr)  
                <type-data>, INTENT(INOUT) :: data  
                INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries  
                INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_id  
                INTEGER(Zoltan_INT), INTENT(IN) :: size  
                INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: buf  
                INTEGER(Zoltan_INT), INTENT(OUT) :: ierr  
  
                <type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),  
                DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x)  
                where x is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.
```

A **ZOLTAN_UNPACK_OBJ_FN** query function allows the application to tell Zoltan how to copy all needed data for a given object from a communication buffer into the application's data structure. This operation is needed as the final step of importing objects during data migration. The query function may also perform other computation, such as building request lists for related data. This routine is called by [Zoltan_Migrate](#) for each object to be received by the processor. (Note: a local ID for the object is not included in this function, as the local ID is local to the exporting, not the importing, processor.)

Function Type: **ZOLTAN_UNPACK_OBJ_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>global_id</i>	The global ID of the object whose data has been received in the communication buffer.
<i>size</i>	The size (in bytes) of the object's data in the communication buffer. This value is at least as large as the value returned by the ZOLTAN_OBJ_SIZE_MULTI_FN or ZOLTAN_OBJ_SIZE_FN query function; it may be slightly larger due to padding for data alignment in the buffer.
<i>buf</i>	The starting address of the communication buffer for this object.
<i>ierr</i>	Error code to be set by function.

C and C++:

typedef void **ZOLTAN_UNPACK_OBJ_MULTI_FN** (
void *data,
int num_gid_entries,
int num_ids,
[ZOLTAN_ID_PTR](#) global_ids,
int *sizes,
int *idx,
char *buf,
int *ierr);

FORTRAN:

SUBROUTINE *Unpack_Obj_Multi*(data, num_gid_entries, num_ids, global_ids, sizes, idx, buf,
ierr)
<type-data>, INTENT(INOUT) :: data
INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries
INTEGER(Zoltan_INT), INTENT(IN) :: num_ids
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: global_ids
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: sizes
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: idx
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: buf
INTEGER(Zoltan_INT), INTENT(OUT) :: ierr

<type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT),
DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x)
where x is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_UNPACK_OBJ_MULTI_FN** query function is the multiple-ID version of a [ZOLTAN_UNPACK_OBJ_FN](#). It allows the application to tell Zoltan how to copy all needed data for a given list of objects from a communication buffer into the application's data structure.

Function Type:	ZOLTAN_UNPACK_OBJ_MULTI_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_ids</i>	The number of objects to be unpacked.
<i>global_ids</i>	An array of global IDs of the objects. The ID for the <i>i</i> -th object begins in <i>global_ids[i*num_gid_entries]</i> .
<i>sizes</i>	An array containing the per-object sizes (in bytes) of the communication buffer for each object. Each value is at least as large as the corresponding value returned by the ZOLTAN_OBJ_SIZE_MULTI_FN or ZOLTAN_OBJ_SIZE_FN query function; it may be slightly larger due to padding for data alignment in the buffer.
<i>idx</i>	For each object, an index into the <i>buf</i> array giving the starting location of that object's data. Data for the <i>i</i> -th object are stored in <i>buf[idx[i]]</i> , <i>buf[idx[i]+1]</i> , ..., <i>buf[idx[i]+sizes[i]-1]</i> . Because Zoltan adds some tag information to packed data, <i>idx[i] != sum[j=0,i-1](sizes[j])</i> .
<i>buf</i>	The address of the communication buffer from which data is unpacked.
<i>ierr</i>	Error code to be set by function.

C and C++:

typedef void **ZOLTAN_PRE_MIGRATE_PP_FN** (
void *data,
int num_gid_entries,

```
int num_lid_entries,  
int num_import,  
ZOLTAN\_ID\_PTR import_global_ids,  
ZOLTAN\_ID\_PTR import_local_ids,  
int *import_procs,  
int *import_to_part,  
int num_export,  
ZOLTAN\_ID\_PTR export_global_ids,  
ZOLTAN\_ID\_PTR export_local_ids,  
int *export_procs,  
int *export_to_part,  
int *ierr);
```

FORTTRAN: SUBROUTINE *Pre_Migrate_PP*(data, num_gid_entries, num_lid_entries, num_import,
import_global_ids, import_local_ids, import_procs, import_to_part, num_export,
export_global_ids, export_local_ids, export_procs, export_to_part, ierr)
 <type-data>, INTENT(INOUT) :: data
 INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries
 INTEGER(Zoltan_INT), INTENT(IN) :: num_import, num_export
 INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_global_ids, export_global_ids
 INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_local_ids, export_local_ids
 INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_procs, export_procs
 INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_to_part, export_to_part
 INTEGER(Zoltan_INT), INTENT(OUT) :: ierr

<type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_PRE_MIGRATE_PP_FN** query function performs any pre-processing desired by the application. If it is registered, it is called at the beginning of the [Zoltan_Migrate](#) routine. The arguments passed to [Zoltan_Migrate](#) are made available for use in the pre-processing routine.

Function Type: **ZOLTAN_PRE_MIGRATE_PP_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES .
<i>num_import</i>	The number of objects that will be received by this processor.
<i>import_global_ids</i>	An array of <i>num_import</i> global IDs of objects to be received by this processor. This array may be NULL, as the processor does not necessarily need to know which objects it will receive.
<i>import_local_ids</i>	An array of <i>num_import</i> local IDs of objects to be received by this processor. This array may be NULL, as the processor does not necessarily need to know which objects it will receive.
<i>import_procs</i>	An array of size <i>num_import</i> listing the processor IDs of the source processors. This array may be NULL, as the processor does not necessarily need to know which objects it will receive.
<i>import_to_part</i>	An array of size <i>num_import</i> listing the parts to which objects will be imported. This array may be NULL, as the processor does not necessarily need to know from which objects it will receive.

<i>num_export</i>	The number of objects that will be sent from this processor to other processors.
<i>export_global_ids</i>	An array of <i>num_export</i> global IDs of objects to be sent from this processor.
<i>export_local_ids</i>	An array of <i>num_export</i> local IDs of objects to be sent from this processor.
<i>export_procs</i>	An array of size <i>num_export</i> listing the processor IDs of the destination processors.
<i>export_to_part</i>	An array of size <i>num_export</i> listing the parts to which objects will be sent.
<i>ierr</i>	Error code to be set by function.

Default: No pre-processing is done if a **ZOLTAN_PRE_MIGRATE_PP_FN** is not registered.

C and C++:	<pre>typedef void ZOLTAN_MID_MIGRATE_PP_FN (void *data, int num_gid_entries, int num_lid_entries, int num_import, ZOLTAN_ID_PTR import_global_ids, ZOLTAN_ID_PTR import_local_ids, int *import_procs, int *import_to_part, int num_export, ZOLTAN_ID_PTR export_global_ids, ZOLTAN_ID_PTR export_local_ids, int *export_procs, int *export_to_part, int *ierr);</pre>
FORTTRAN:	<pre>SUBROUTINE Mid_Migrate_PP(data, num_gid_entries, num_lid_entries, num_import, import_global_ids, import_local_ids, import_procs, import_to_part, num_export, export_global_ids, export_local_ids, export_procs, export_to_part, ierr) <type-data>, INTENT(INOUT) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries INTEGER(Zoltan_INT), INTENT(IN) :: num_import, num_export INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_global_ids, export_global_ids INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_local_ids, export_local_ids INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_procs, export_procs INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_to_part, export_to_part INTEGER(Zoltan_INT), INTENT(OUT) :: ierr</pre> <p><type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where <i>x</i> is 1, 2, 3 or 4. See the section on Fortran query functions for an explanation.</p>

A **ZOLTAN_MID_MIGRATE_PP_FN** query function performs any processing desired by the application between the packing and unpacking of objects being migrated. If it is registered, it is called after export objects are packed in [Zoltan_Migrate](#); imported objects are unpacked after the **ZOLTAN_MID_MIGRATE_PP_FN** query function is called. The arguments passed to [Zoltan_Migrate](#) are made available for use in the processing routine.

Function Type:	ZOLTAN_MID_MIGRATE_PP_FN_TYPE
Arguments:	
<i>data</i>	Pointer to user-defined data.

<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES .
<i>num_import</i>	The number of objects that will be received by this processor.
<i>import_global_ids</i>	An array of <i>num_import</i> global IDs of objects to be received by this processor. This array may be NULL, as the processor does not necessarily need to know which objects it will receive.
<i>import_local_ids</i>	An array of <i>num_import</i> local IDs of objects to be received by this processor. This array may be NULL, as the processor does not necessarily need to know which objects it will receive.
<i>import_procs</i>	An array of size <i>num_import</i> listing the processor IDs of the source processors. This array may be NULL, as the processor does not necessarily need to know which objects it will receive.
<i>import_to_part</i>	An array of size <i>num_import</i> listing the parts to which objects will be imported. This array may be NULL, as the processor does not necessarily need to know from which objects it will receive.
<i>num_export</i>	The number of objects that will be sent from this processor to other processors.
<i>export_global_ids</i>	An array of <i>num_export</i> global IDs of objects to be sent from this processor.
<i>export_local_ids</i>	An array of <i>num_export</i> local IDs of objects to be sent from this processor.
<i>export_procs</i>	An array of size <i>num_export</i> listing the processor IDs of the destination processors.
<i>export_to_part</i>	An array of size <i>num_export</i> listing the parts to which objects will be sent.
<i>ierr</i>	Error code to be set by function.

Default:

No processing is done if a **ZOLTAN_MID_MIGRATE_PP_FN** is not registered.

C and C++:	<pre>typedef void ZOLTAN_POST_MIGRATE_PP_FN (void *data, int num_gid_entries, int num_lid_entries, int num_import, ZOLTAN_ID_PTR import_global_ids, ZOLTAN_ID_PTR import_local_ids, int *import_procs, int *import_to_part, int num_export, ZOLTAN_ID_PTR export_global_ids, ZOLTAN_ID_PTR export_local_ids, int *export_procs, int *export_to_part, int *ierr);</pre>
FORTTRAN:	<pre>SUBROUTINE Post_Migrate_PP(data, num_gid_entries, num_lid_entries, num_import, import_global_ids, import_local_ids, import_procs, import_to_part, num_export, export_global_ids, export_local_ids, export_procs, export_to_part, ierr) <type-data>, INTENT(INOUT) :: data INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries INTEGER(Zoltan_INT), INTENT(IN) :: num_import, num_export INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_global_ids, export_global_ids INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_local_ids, export_local_ids INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_procs, export_procs</pre>

INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_to_part, export_to_part
INTEGER(Zoltan_INT), INTENT(OUT) :: ierr

<type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_POST_MIGRATE_PP_FN** query function performs any post-processing desired by the application. If it is registered, it is called at the end of the [Zoltan_Migrate](#) routine. The arguments passed to [Zoltan_Migrate](#) are made available for use in the post-processing routine.

Function Type: **ZOLTAN_POST_MIGRATE_PP_FN_TYPE**

Arguments:

<i>data</i>	Pointer to user-defined data.
<i>num_gid_entries</i>	The number of array entries used to describe a single global ID. This value is the maximum value over all processors of the parameter NUM_GID_ENTRIES .
<i>num_lid_entries</i>	The number of array entries used to describe a single local ID. This value is the maximum value over all processors of the parameter NUM_LID_ENTRIES .
<i>num_import</i>	The number of objects that will be received by this processor.
<i>import_global_ids</i>	An array of <i>num_import</i> global IDs of objects to be received by this processor. This array may be NULL, as the processor does not necessarily need to know which objects it will receive.
<i>import_local_ids</i>	An array of <i>num_import</i> local IDs of objects to be received by this processor. This array may be NULL, as the processor does not necessarily need to know which objects it will receive.
<i>import_procs</i>	An array of size <i>num_import</i> listing the processor IDs of the source processors. This array may be NULL, as the processor does not necessarily need to know which objects it will receive.
<i>import_to_part</i>	An array of size <i>num_import</i> listing the parts to which objects will be imported. This array may be NULL, as the processor does not necessarily need to know from which objects it will receive.
<i>num_export</i>	The number of objects that will be sent from this processor to other processors.
<i>export_global_ids</i>	An array of <i>num_export</i> global IDs of objects to be sent from this processor.
<i>export_local_ids</i>	An array of <i>num_export</i> local IDs of objects to be sent from this processor.
<i>export_procs</i>	An array of size <i>num_export</i> listing the processor IDs of the destination processors.
<i>export_to_part</i>	An array of size <i>num_export</i> listing the parts to which objects will be sent.
<i>ierr</i>	Error code to be set by function.

Default:

No post-processing is done if a **ZOLTAN_POST_MIGRATE_PP_FN** is not registered.

C: typedef void **ZOLTAN_PRE_MIGRATE_FN** (
 void *data,
 int num_gid_entries,
 int num_lid_entries,
 int num_import,
 [ZOLTAN_ID_PTR](#) import_global_ids,
 [ZOLTAN_ID_PTR](#) import_local_ids,
 int *import_procs,
 int num_export,

```
ZOLTAN\_ID\_PTR export_global_ids,  
ZOLTAN\_ID\_PTR export_local_ids,  
int *export_procs,  
int *ierr);
```

FORTTRAN: SUBROUTINE *Pre_Migrate*(data, num_gid_entries, num_lid_entries, num_import, import_global_ids, import_local_ids, import_procs, num_export, export_global_ids, export_local_ids, export_procs, ierr)
<type-data>, INTENT(INOUT) :: data
INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries
INTEGER(Zoltan_INT), INTENT(IN) :: num_import, num_export
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_global_ids, export_global_ids
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_local_ids, export_local_ids
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_procs, export_procs
INTEGER(Zoltan_INT), INTENT(OUT) :: ierr

<type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_PRE_MIGRATE_FN** query function performs any pre-processing desired by applications using [Zoltan Help Migrate](#). Its function is analogous to [ZOLTAN_PRE_MIGRATE_PP_FN](#), but it cannot be used with [Zoltan Migrate](#).

Function Type: ZOLTAN_PRE_MIGRATE_FN_TYPE
Arguments:

All arguments are analogous to those in [ZOLTAN_PRE_MIGRATE_PP_FN](#). Part-assignment arguments *import_to_part* and *export_to_part* are not included, as processor and parts numbers are considered to be the same in [Zoltan Help Migrate](#).

Default:
No pre-processing is done if a **ZOLTAN_PRE_MIGRATE_FN** is not registered.

C: typedef void **ZOLTAN_MID_MIGRATE_FN** (
void *data,
int num_gid_entries,
int num_lid_entries,
int num_import,
[ZOLTAN_ID_PTR](#) import_global_ids,
[ZOLTAN_ID_PTR](#) import_local_ids,
int *import_procs,
int num_export,
[ZOLTAN_ID_PTR](#) export_global_ids,
[ZOLTAN_ID_PTR](#) export_local_ids,
int *export_procs,
int *ierr);

FORTTRAN: SUBROUTINE *Mid_Migrate*(data, num_gid_entries, num_lid_entries, num_import, import_global_ids, import_local_ids, import_procs, num_export, export_global_ids, export_local_ids, export_procs, ierr)
<type-data>, INTENT(INOUT) :: data
INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries
INTEGER(Zoltan_INT), INTENT(IN) :: num_import, num_export

INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_global_ids, export_global_ids
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_local_ids, export_local_ids
INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_procs, export_procs
INTEGER(Zoltan_INT), INTENT(OUT) :: ierr

<type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_MID_MIGRATE_FN** query function performs any mid-migration processing desired by applications using [Zoltan Help Migrate](#). Its function is analogous to [ZOLTAN_MID_MIGRATE_PP_FN](#), but it cannot be used with [Zoltan Migrate](#).

Function Type: **ZOLTAN_MID_MIGRATE_FN_TYPE**

Arguments: All arguments are analogous to those in [ZOLTAN_MID_MIGRATE_PP_FN](#). Part-assignment arguments *import_to_part* and *export_to_part* are not included, as processor and parts numbers are considered to be the same in [Zoltan Help Migrate](#).

Default: No processing is done if a **ZOLTAN_MID_MIGRATE_FN** is not registered.

C: typedef void **ZOLTAN_POST_MIGRATE_FN** (
 void *data,
 int num_gid_entries,
 int num_lid_entries,
 int num_import,
 [ZOLTAN_ID_PTR](#) import_global_ids,
 [ZOLTAN_ID_PTR](#) import_local_ids,
 int *import_procs,
 int num_export,
 [ZOLTAN_ID_PTR](#) export_global_ids,
 [ZOLTAN_ID_PTR](#) export_local_ids,
 int *export_procs,
 int *ierr);

FORTTRAN: SUBROUTINE **Post_Migrate**(data, num_gid_entries, num_lid_entries, num_import, import_global_ids, import_local_ids, import_procs, num_export, export_global_ids, export_local_ids, export_procs, ierr)
 <type-data>, INTENT(INOUT) :: data
 INTEGER(Zoltan_INT), INTENT(IN) :: num_gid_entries, num_lid_entries
 INTEGER(Zoltan_INT), INTENT(IN) :: num_import, num_export
 INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_global_ids, export_global_ids
 INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_local_ids, export_local_ids
 INTEGER(Zoltan_INT), INTENT(IN), DIMENSION(*) :: import_procs, export_procs
 INTEGER(Zoltan_INT), INTENT(OUT) :: ierr

<type-data> can be any of INTEGER(Zoltan_INT), DIMENSION(*) or REAL(Zoltan_FLOAT), DIMENSION(*) or REAL(Zoltan_DOUBLE), DIMENSION(*) or TYPE(Zoltan_User_Data_x) where *x* is 1, 2, 3 or 4. See the section on [Fortran query functions](#) for an explanation.

A **ZOLTAN_POST_MIGRATE_FN** query function performs any post-processing desired by applications using

[Zoltan_Help_Migrate](#). Its function is analogous to [ZOLTAN_POST_MIGRATE_PP_FN](#), but it cannot be used with [Zoltan_Migrate](#).

Function Type: ZOLTAN_POST_MIGRATE_FN_TYPE

Arguments: All arguments are analogous to those in [ZOLTAN_POST_MIGRATE_PP_FN](#). Part-assignment arguments *import_to_part* and *export_to_part* are not included, as processor and parts numbers are considered to be the same in [Zoltan_Help_Migrate](#).

Default: No post-processing is done if a [ZOLTAN_POST_MIGRATE_FN](#) is not registered.

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Zoltan Parameters and Output Levels

The behavior of Zoltan is controlled by several [parameters](#) and [debugging-output levels](#). These parameters can be set by calls to [Zoltan_Set_Param](#). Reasonable [default values](#) for all parameters are specified by Zoltan. Many of the parameters are specific to individual algorithms, and are listed in the descriptions of those algorithms. However, the parameters below have meaning across the entire library.

General Parameters

The following parameters apply to the entire Zoltan library. While reasonable [default values](#) for all parameters are specified by Zoltan, applications can change these values through calls to [Zoltan_Set_Param](#).

Parameters:

<i>NUM_GID_ENTRIES</i>	The number of unsigned integers that should be used to represent a global identifier (ID). Values greater than zero are accepted.
<i>NUM_LID_ENTRIES</i>	The number of unsigned integers that should be used to represent a local identifier (ID). Values greater than or equal to zero are accepted.
<i>DEBUG_LEVEL</i>	An integer indicating how much debugging information is printed by Zoltan. Higher values of <i>DEBUG_LEVEL</i> produce more output and potentially slow down Zoltan's computations. The least output is produced when <i>DEBUG_LEVEL</i> = 0. <i>DEBUG_LEVEL</i> primarily controls Zoltan's behavior; most algorithms have their own parameters to control their output level. Values used within Zoltan are listed below . Note: Because some debugging levels use processor synchronization, all processors should use the same value of <i>DEBUG_LEVEL</i> . Processor number from which trace output should be printed when <i>DEBUG_LEVEL</i> is 5.
<i>DEBUG_PROCESSOR</i>	
<i>DEBUG_MEMORY</i>	Integer indicating the amount of low-level debugging information about memory-allocation should be kept by Zoltan's Memory Management utilities . Valid values are 0, 1, 2, and 3.
<i>OBJ_WEIGHT_DIM</i>	The number of weights (to be supplied by the user in a query function) associated with an object. If this parameter is zero, all objects have equal weight. Some algorithms may not support multiple (multidimensional) weights.
<i>EDGE_WEIGHT_DIM</i>	The number of weights associated with an edge. If this parameter is zero, all edges have equal weight. Many algorithms do not support multiple (multidimensional) weights.
<i>TIMER</i>	The timer with which you wish to measure time. Valid choices are <i>wall</i> (based on <i>MPI_Wtime</i>), <i>cpu</i> (based on the ANSI C library function <i>clock</i>), and <i>user</i> . The resolution may be poor, as low as 1/60th of a second, depending upon your platform.

Default Values:

[*NUM_GID_ENTRIES*](#) = 1
[*NUM_LID_ENTRIES*](#) = 1
[*DEBUG_LEVEL*](#) = 1
[*DEBUG_PROCESSOR*](#) = 0
[*DEBUG_MEMORY*](#) = 1
[*OBJ_WEIGHT_DIM*](#) = 0
[*EDGE_WEIGHT_DIM*](#) = 0
[*TIMER*](#) = wall

Debugging Levels in Zoltan

The `DEBUG_LEVEL` parameter determines how much debugging information is printed to `stdout` by Zoltan. It is set by a call to [Zoltan_Set_Param](#). Higher values of `DEBUG_LEVEL` produce more output and can slow down Zoltan's computations, especially when the output is printed by one processor at a time. The least output is produced when `DEBUG_LEVEL = 0`.

Descriptions of the output produced by Zoltan for each value of `DEBUG_LEVEL` are included below. For a given `DEBUG_LEVEL` value n , all output for values less than or equal to n is produced.

Some high debugging levels use processor synchronization to force processors to write one-at-a-time. For example, when `DEBUG_LEVEL` is greater than or equal to eight, each processor writes its list in turn so that the lists from all processors are not jumbled together in the output. This synchronization requires all processors to use the same value of `DEBUG_LEVEL`.

`DEBUG_LEVEL` Output Produced

- 0 Quiet mode; no output unless an error or warning is produced.
- 1 Values of all parameters set by [Zoltan_Set_Param](#) and used by Zoltan.
- 2 Timing information for Zoltan's main routines.
- 3 Timing information within Zoltan's algorithms (support by algorithms is optional).
- 4
- 5 Trace information (enter/exit) for major Zoltan interface routines (printed by the processor specified by the [DEBUG_PROCESSOR](#) parameter).
- 6 Trace information (enter/exit) for major Zoltan interface routines (printed by all processors).
- 7 More detailed trace information in major Zoltan interface routines.
- 8 List of objects to be imported to and exported from each processor. ¹
- 9
- 10 Maximum debug output; may include algorithm-specific output. ¹

¹ *Output may be serialized; that is, one processor may have to complete its output before the next processor is allowed to begin its output. This serialization is not scalable and can significantly increase execution time on large number of processors.*

Load-Balancing Algorithms and Parameters

The following dynamic load-balancing algorithms are currently included in the Zoltan library:

[Simple Partitioners for Testing](#)

- [Block Partitioning \(BLOCK\)](#)
- [Cyclic Partitioning \(CYCLIC\)](#)
- [Random Partitioning \(RANDOM\)](#)

[Geometric \(Coordinate-based\) Partitioners](#)

- [Recursive Coordinate Bisection \(RCB\)](#)
- [Recursive Inertial Bisection \(RIB\)](#)
- [Hilbert Space-Filling Curve Partitioning \(HSFC\)](#)
- [Refinement Tree Based Partitioning \(REFTREE\)](#)

[Hypergraph Partitioning, Repartitioning and Refinement \(HYPERGRAPH\)](#)

- [PHG](#)
- [PaToH](#)

[Graph Partitioning and Repartitioning \(GRAPH\)](#)

- [PHG](#)
- [ParMETIS](#)
- [Scotch](#)

[Hybrid Hierarchical Partitioning \(HIER\)](#)

The parenthetical string is the parameter value for [LB_METHOD](#) parameter; the parameter is set through a call to [Zoltan_Set_Param](#).

For further analysis and discussion of some of the algorithms, see [[Hendrickson and Devine](#)].

Load-Balancing Parameters

While the overall behavior of Zoltan is controlled by [general Zoltan parameters](#), the behavior of each load-balancing method is controlled by parameters specific to partitioning which are also set by calls to [Zoltan_Set_Param](#). Many of these parameters are specific to individual partitioning algorithms, and are listed in the descriptions of the individual algorithms. However, several have meaning across multiple partitioning algorithms. These load-balancing parameters are described below. Unless indicated otherwise, these parameters apply to both [Zoltan_LB_Partition](#) and [Zoltan_LB_Balance](#).

Parameters:

<code>LB_METHOD</code>	The load-balancing algorithm used by Zoltan is specified by this parameter. Valid values are
	BLOCK (for block partitioning),

RANDOM (for [random partitioning](#)),
RCB (for [recursive coordinate bisection](#)),
RIB (for [recursive inertial bisection](#)),
HSFC (for [Hilbert space-filling curve partitioning](#)),
REFTREE (for [refinement tree based partitioning](#))
GRAPH (to choose from collection of methods for [graphs](#)),
HYPERGRAPH (to choose from a collection of methods for [hypergraphs](#)),
HIER (for hybrid [hierarchical partitioning](#))
NONE (for no load balancing).

LB_APPROACH

The desired load balancing approach. Only *LB_METHOD* = [HYPERGRAPH](#) or [GRAPH](#) uses the *LB_APPROACH* parameter. Valid values are

PARTITION (Partition "from scratch," not taking into account the current data distribution; this option is recommended for static load balancing.)
REPARTITION (Partition but take into account current data distribution to keep data migration low; this option is recommended for dynamic load balancing.)
REFINE (Quickly improve the current data distribution.)

NUM_GLOBAL_PARTS

The total number of parts to be generated by a call to [Zoltan_LB_Partition](#). Integer values greater than zero are accepted. Not valid for [Zoltan_LB_Balance](#).

NUM_LOCAL_PARTS

The number of parts to be generated on this processor by a call to [Zoltan_LB_Partition](#). Integer values greater than or equal to zero are accepted. Not valid for [Zoltan_LB_Balance](#). If any processor sets this parameter, NUM_LOCAL_PARTS is assumed to be zero on processors not setting this parameter.

RETURN_LISTS

The lists returned by calls to [Zoltan_LB_Partition](#) or [Zoltan_LB_Balance](#). Valid values are

"IMPORT", to return only information about objects to be imported to a processor
"EXPORT", to return only information about objects to be exported from a processor
"ALL", or "IMPORT AND EXPORT" (or any string with both "IMPORT" and "EXPORT" in it) to return both import and export information
"PARTS" (or "PART ASSIGNMENT" or any string with "PART" in it) to return the new process and part assignment of every local object, including those not being exported.
"NONE", to return neither import nor export information

REMAP

Within [Zoltan_LB_Partition](#) or [Zoltan_LB_Balance](#), renumber parts to maximize overlap between the old decomposition and the new decomposition (to reduce data movement from old to new decompositions). Valid values are "0" (no remapping) or "1" (remapping). Part assignments from [ZOLTAN_PART_MULTIFN](#) or [ZOLTAN_PART_FN](#) query functions can be used in remapping if provided; otherwise, processor numbers are used as part numbers. Requests for remapping are ignored when, in the new decomposition,

a part is spread across multiple processors or part sizes are specified using [Zoltan LB Set Part Sizes](#).

IMBALANCE_TOL

The amount of load imbalance the partitioning algorithm should deem acceptable. The load on each processor is computed as the sum of the weights of objects it is assigned. The imbalance is then computed as the maximum load divided by the average load. An value for *IMBALANCE_TOL* of 1.2 indicates that 20% imbalance is OK; that is, the maximum over the average shouldn't exceed 1.2.

MIGRATE_ONLY_PROC_CHANGES

If this value is set to TRUE (non-zero), Zoltan's migration functions will migrate only objects moving to new processors. They will not migrate objects for which only the part number has changed; the objects' processor numbers must change as well. If this value is set to FALSE (zero), Zoltan's migration functions will migrate all objects with new part or processor assignments.

AUTO_MIGRATE

If this value is set to TRUE (non-zero), Zoltan will automatically perform the data migration during calls to [Zoltan LB Partition](#) or [Zoltan LB Balance](#). A full discussion of automatic migration can be found in the description of the [migration interface functions](#).

Default Values:

LB_METHOD = RCB

LB_APPROACH = REPARTITION

NUM_GLOBAL_PARTS = Number of processors specified in [Zoltan Create](#).

NUM_LOCAL_PARTS = 1

RETURN_LISTS = ALL

REMAP = 1

IMBALANCE_TOL = 1.1

MIGRATE_ONLY_PROC_CHANGES = 1

AUTO_MIGRATE = FALSE

Simple Partitioners for Testing

Zoltan includes two very simple partitioners for testing initial implementations of Zoltan in applications. These partitioners are intended only for testing and to serve as examples. They use neither geometry nor connectivity (graph/hypergraph), so they require very few query functions (only two!).

- [Block partitioning](#) (BLOCK)
- [Random partitioning](#) (RANDOM)

[[Table of Contents](#) | [Next: Block Partitioning](#) | [Previous: Load-Balancing Algorithms and Parameters](#) | [Privacy and Security](#)]

Block

A simple partitioner based on block partitioning of the objects. It is mainly intended for testing. It uses neither geometry nor connectivity (graph/hypergraph), so it requires very few query functions. The block strategy is as follows: Consider all objects (on all processors) as a linear sequence. Assign the first block of $n/\text{num_parts}$ objects to the first part, the next block to the second, and so on. Block is smart enough to generalize this method to handle vertex weights and target part sizes. Only a single weight per object ($\text{Obj_Weight_Dim}=1$) is currently supported.

Method String:	Block
Parameters:	
Required Query Functions:	ZOLTAN_NUM_OBJ_FN ZOLTAN_OBJ_LIST_FN

Cyclic

A simple partitioner based on cyclic (round robin) partitioning of the objects. It uses neither geometry nor connectivity (graph/hypergraph), so it requires very few query functions. This method currently does not take into account target part sizes nor weights, so the parts may not be balanced in this case!

Note that the partitioning is cyclic with respect to the local order of objects, and NOT by the global ids.

Method String:	Cyclic
Parameters:	
Required Query Functions:	ZOLTAN_NUM_OBJ_FN ZOLTAN_OBJ_LIST_FN

Random

Random is not really a load-balancing algorithm, and should be used only for testing! It takes each object and randomly assigns it to a new part. Via a parameter, one can alternatively choose to randomly perturb only a fraction of the objects. The random method does not use weights and does not attempt to achieve load balance.

Method String:	Random
Parameters:	
<i>RANDOM_MOVE_FRACTION</i>	The fraction of objects to randomly move. 1.0 = move all; 0.0 = move nothing
Default:	<i>RANDOM_MOVE_FRACTION</i> = 1.0
Required Query Functions:	ZOLTAN_NUM_OBJ_FN ZOLTAN_OBJ_LIST_FN

Geometric (Coordinate-based) Partitioners

Geometric partitioners divide data into parts based on the physical coordinates of the data. Objects assigned to a single part tend to be physically close to each other in space. Such partitioners are very useful for applications that don't have explicit connectivity information (such as particle methods) or for which geometric locality is important (such as contact detection). They are also widely used in adaptive finite element methods because, in general, they execute very quickly and yield moderately good partition quality.

The geometric methods are the easiest non-trivial partitioners to incorporate into applications, as they require only four callbacks: two returning [object information](#) and two returning [coordinate information](#).

We group [refinement-tree partitioning](#) for adaptive mesh refinement applications into the geometric partitioners because it uses geometric information to determine an initial ordering for coarse elements of adaptively refined meshes. The refinement-tree partitioner also requires [tree-based callbacks](#) with connectivity information between coarse and fine elements in refined meshes.

- [Recursive Coordinate Bisection](#) (RCB)
- [Recursive Inertial Bisection](#) (RIB)
- [Hilbert Space-Filling Curve Partitioning](#) (HSFC)
- [Refinement Tree Based Partitioning](#) (Reftree)

Recursive Coordinate Bisection (RCB)

An implementation of Recursive Coordinate Bisection (RCB) due to Steve Plimpton of Sandia National Laboratories is included in Zoltan. RCB was first proposed as a static load-balancing algorithm by [Berger and Bokhari](#), but is attractive as a dynamic load-balancing algorithm because it implicitly produces incremental partitions. In RCB, the computational domain is first divided into two regions by a cutting plane orthogonal to one of the coordinate axes so that half the work load is in each of the sub-regions. The splitting direction is determined by computing in which coordinate direction the set of objects is most elongated, based upon the geometric locations of the objects. The sub-regions are then further divided by recursive application of the same splitting algorithm until the number of sub-regions equals the number of processors. Although this algorithm was first devised to cut into a number of sets which is a power of two, the set sizes in a particular cut needn't be equal. By adjusting the part sizes appropriately, any number of equally-sized sets can be created. If the parallel machine has processors with different speeds, sets with nonuniform sizes can also be easily generated. The Zoltan implementation of RCB has several parameters which can be modified by the [Zoltan Set Param](#) function. A recent feature is that RCB allows multiple weights; that is, one can balance with respect to several load criteria simultaneously. Note that there is no guarantee that a desired load balance tolerance can be achieved using RCB, especially in the multiconstraint case.

Information about the sub-regions generated by RCB can be obtained by an application through calls to [Zoltan_RCB_Box](#). This function is not required to perform load balancing; it only provides auxiliary information to an application.

Method String:	RCB
Parameters:	
<i>RCB_OVERALLOC</i>	The amount by which to over-allocate temporary storage arrays for objects within the RCB algorithm when additional storage is due to changes in processor assignments. 1.0 = no extra storage allocated; 1.5 = 50% extra storage; etc.
<i>RCB_REUSE</i>	Flag to indicate whether to use previous cuts as initial guesses for the current RCB invocation. 0 = don't use previous cuts; 1 = use previous cuts.
<i>RCB_OUTPUT_LEVEL</i>	Flag controlling the amount of timing and diagnostic output the routine produces. 0 = no output; 1 = print summary; 2 = print data for each processor.
<i>CHECK_GEOM</i>	Flag controlling the invocation of input and output error checking. 0 = don't do checking; 1 = do checking.
<i>KEEP_CUTS</i>	Should information about the cuts determining the RCB decomposition be retained? It costs a bit of time to do so, but this information is necessary if application wants to add more objects to the decomposition via calls to Zoltan_LB_Point_PP_Assign or to Zoltan_LB_Box_PP_Assign . 0 = don't keep cuts; 1 = keep cuts.
<i>AVERAGE_CUTS</i>	When set to one, coordinates of RCB cutting planes are computed to be the average of the coordinates of the closest object on each side of the cut. Otherwise, coordinates of cutting planes may equal those of one of the closest objects. 0 = don't average cuts; 1 = average cuts.
<i>RCB_LOCK_DIRECTIONS</i>	Flag that determines whether the order of the directions of the cuts is kept constant after they are determined the first time RCB is called. 0 = don't lock directions; 1 = lock directions.
<i>RCB_SET_DIRECTIONS</i>	If this flag is set, the order of cuts is changed so that all of the cuts in any direction

are done as a group. The number of cuts in each direction is determined and then the value of the parameter is used to determine the order that those cuts are made in. When 1D and 2D problems are partitioned, the directions corresponding to unused dimensions are ignored.

0 = don't order cuts; 1 = xyz; 2 = xzy; 3 = yzx; 4 = yxz; 5 = zxy; 6 = zyx;

Flag controlling the shape of the resulting regions. If this option is specified, then when a cut is made, all of the dots located on the cut are moved to the same side of the cut. The resulting regions are then rectilinear. When these dots are treated as a group, then the resulting load balance may not be as good as when the group of dots is split by the cut.

0 = move dots individually; 1 = move dots in groups.

When a 3 dimensional geometry is almost flat, it may make more sense to treat it as a 2 dimensional geometry when applying the RCB algorithm. In this case, a 2 dimensional RCB calculation is applied to a plane that corresponds with the geometry. (This results in cuts that, while still orthogonal, may no longer be axis aligned.) If this parameter is set to **1**, a 3 dimensional geometry will be treated as 2 dimensional if it is very flat, or 1 dimensional if it is very thin. A 2 dimensional geometry will be treated as 1 dimensional if it is very thin.

If the **REDUCE_DIMENSIONS** parameter is set, then this parameter determines when a geometry is considered to be degenerate. A bounding box which is oriented to the geometry is constructed, and the lengths of its sides are tested against a ratio of 1 : **DEGENERATE_RATIO**.

Flag indicating whether the bounding box of set of parts is recomputed at each level of recursion. By default, the longest direction of the bounding box is cut during bisection. Recomputing the bounding box at each level of recursion can produce more effective cut directions for unusually shaped geometries; the computation does, however, take additional time and communication, and may cause cut directions to vary from one invocation of RCB to the next.

0 = don't recompute the bounding box; 1 = recompute the box.

In the multiconstraint case, are the object weights comparable? Do they have the same units and is the scaling meaningful? For example, if the jth weight corresponds to the expected time in phase j (measured in seconds), set this parameter to 1. (0 = incomparable, 1 = comparable)

Norm used in multicriteria algorithm; this determines how to balance the different weight constraints. Valid values are 1,2, and 3. Roughly, if the weights correspond to different phases, then the value 1 (1-norm) tries to minimize the total time (sum over all phases) while the value 3 (max-norm) attempts to minimize the worst imbalance in any phase. The 2-norm does something in between. Try a different value if you're not happy with the balance.

Maximum allowed ratio between the largest and smallest side of a subdomain. Must be > 1.

RCB_OVERALLOC = 1.2

RCB_REUSE = 0

RCB_OUTPUT_LEVEL = 0

CHECK_GEOM = 1

KEEP_CUTS = 0

AVERAGE_CUTS = 0

RCB_LOCK_DIRECTIONS = 0

REDUCE_DIMENSIONS = 0

DEGENERATE_RATIO = 1.0

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

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RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

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OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

OBJ_WEIGHTS_COMPARABLE = 0

RCB_MULTICRITERIA_NORM = 1

RCB_MAX_ASPECT_RATIO = 1.0

RCB_RECOMPUTE_BOX = 0

RCB_RECTILINEAR_BLOCKS = 0

DEGENERATE_RATIO = 10
RCB_SET_DIRECTIONS = 0
RCB_RECTILINEAR_BLOCKS = 0
RCB_RECOMPUTE_BOX = 0
OBJ_WEIGHTS_COMPARABLE = 0
RCB_MULTICRITERIA_NORM = 1
RCB_MAX_ASPECT_RATIO = 10

Required Query Functions:

[ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)
[ZOLTAN_NUM_GEOM_FN](#)
[ZOLTAN_GEOM_MULTI_FN](#) or [ZOLTAN_GEOM_FN](#)

```
C:      int Zoltan_RCB_Box (  
        struct Zoltan_Struct * zz,  
        int part,  
        int *ndim,  
        double *xmin,  
        double *ymin,  
        double *zmin,  
        double *xmax,  
        double *ymax,  
        double *zmax);  
  
FORTRAN:  FUNCTION Zoltan_RCB_Box(zz, part, ndim, xmin, ymin, zmin, xmax, ymax, zmax)  
          INTEGER(Zoltan_INT) :: Zoltan_RCB_Box  
          TYPE(Zoltan_Struct), INTENT(IN) :: zz  
          INTEGER(Zoltan_INT), INTENT(IN) :: part  
          INTEGER(Zoltan_INT), INTENT(OUT) :: ndim  
          REAL(Zoltan_DOUBLE), INTENT(OUT) :: xmin, ymin, zmin, xmax, ymax, zmax
```

In many settings, it is useful to know a part's bounding box generated by RCB. This bounding box describes the region of space assigned to a given part. Given an RCB decomposition of space and a part number, **Zoltan_RCB_Box** returns the lower and upper corners of the region of space assigned to the part. To use this routine, the parameter **KEEP_CUTS** must be set to TRUE when the decomposition is generated. This parameter will cause the sequence of geometric cuts to be saved, which is necessary for **Zoltan_RCB_Box** to do its job.

Arguments:

<i>zz</i>	Pointer to the Zoltan structure created by Zoltan_Create .
<i>part</i>	Part number of part for which the bounding box should be returned.
<i>ndim</i>	Upon return, the number of dimensions in the partitioned geometry.
<i>xmin, ymin, zmin</i>	Upon return, the coordinates of the lower extent of bounding box for the part. If the geometry is two-dimensional, <i>zmin</i> is -DBL_MAX. If the geometry is one-dimensional, <i>ymin</i> is -DBL_MAX.
<i>xmax, ymax, zmax</i>	Upon return, the coordinates of the upper extent of bounding box for the part. If the geometry is two-dimensional, <i>zmax</i> is DBL_MAX. If the geometry is one-dimensional, <i>ymax</i> is DBL_MAX.

Returned Value:

int [Error code.](#)

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Recursive Inertial Bisection (RIB)

An implementation of Recursive Inertial Bisection (RIB) is included in Zoltan. RIB was proposed as a load-balancing algorithm by [Williams](#) and later studied by [Taylor and Nour-Omid](#), but its origin is unclear. RIB is similar to RCB in that it divides the domain based on the location of the objects being partitioned by use of cutting planes. In RIB, the computational domain is first divided into two regions by a cutting plane orthogonal to the longest direction of the domain so that half the work load is in each of the sub-regions. The sub-regions are then further divided by recursive application of the same splitting algorithm until the number of sub-regions equals the number of processors. Although this algorithm was first devised to cut into a number of sets which is a power of two, the set sizes in a particular cut needn't be equal. By adjusting the part sizes appropriately, any number of equally-sized sets can be created. If the parallel machine has processors with different speeds, sets with nonuniform sizes can also be easily generated. The Zoltan implementation of RIB has several parameters which can be modified by the [Zoltan_Set_Param](#) function.

Method String: RIB

Parameters:

<i>RIB_OVERALLOC</i>	The amount by which to over-allocate temporary storage arrays for objects within the RIB algorithm when additional storage is due to changes in processor assignments. 1.0 = no extra storage allocated; 1.5 = 50% extra storage; etc.
<i>RIB_OUTPUT_LEVEL</i>	Flag controlling the amount of timing and diagnostic output the routine produces. 0 = no output; 1 = print summary; 2 = print data for each processor.
<i>CHECK_GEOM</i>	Flag controlling the invocation of input and output error checking. 0 = don't do checking; 1 = do checking.
<i>KEEP_CUTS</i>	Should information about the cuts determining the RIB decomposition be retained? It costs a bit of time to do so, but this information is necessary if application wants to add more objects to the decomposition via calls to Zoltan_LB_Point_PP_Assign or to Zoltan_LB_Box_PP_Assign . 0 = don't keep cuts; 1 = keep cuts.
<i>AVERAGE_CUTS</i>	When set to one, coordinates of RIB cutting planes are computed to be the average of the coordinates of the closest object on each side of the cut. Otherwise, coordinates of cutting planes may equal those of one of the closest objects. 0 = don't average cuts; 1 = average cuts.
<i>REDUCE_DIMENSIONS</i>	When a 3 dimensional geometry is almost flat, it may make more sense to treat it as a 2 dimensional geometry when applying the RIB algorithm. (Coordinate values in the omitted direction are ignored for the purposes of partitioning.) If this parameter is set to 1 , a 3 dimensional geometry will be treated as 2 dimensional if it is very flat, or 1 dimensional if it is very thin. A 2 dimensional geometry will be treated as 1 dimensional if it is very thin.
<i>DEGENERATE_RATIO</i>	If the REDUCE_DIMENSIONS parameter is set, then this parameter determines when a geometry is considered to be degenerate. A bounding box which is oriented to the geometry is constructed, and the lengths of its sides are tested against a ratio of 1 : DEGENERATE_RATIO .

Default:

RIB_OVERALLOC = 1.2
RIB_OUTPUT_LEVEL = 0
CHECK_GEOM = 1
KEEP_CUTS = 0
AVERAGE_CUTS = 0

REDUCE_DIMENSIONS = 0
DEGENERATE_RATIO = 10

**Required Query
Functions:**

[ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)
[ZOLTAN_NUM_GEOM_FN](#)
[ZOLTAN_GEOM_MULTI_FN](#) or [ZOLTAN_GEOM_FN](#)

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Hilbert Space Filling Curve (HSFC)

The Inverse Hilbert Space-Filling Curve functions map a point in one, two or three dimensions into the interval [0,1]. The Hilbert functions that map [0, 1] to normal spatial coordinates are also provided. (The one-dimensional inverse Hilbert curve is defined here as the identity function, $f(x)=x$ for all x .)

The HSFC partitioning algorithm seeks to divide [0,1] into P intervals each containing the same weight of objects associated to these intervals by their inverse Hilbert coordinates. N bins are created (where $N > P$) to partition [0,1]. The weights in each bin are summed across all processors. A greedy algorithm sums the bins (from left to right) placing a cut when the desired weight for current part interval is achieved. This process is repeated as needed to improve partitioning tolerance by a technique that maintains the same total number of bins but refines the bins previously containing a cut.

HSFC returns an warning if the final imbalance exceeds the user specified tolerance.

This code implements both the point assign and box assign functionality. The point assign determines an appropriate part (associated with a specific group of processors) for a new point. The box assign determines the list of processors whose associated subdomains intersect the given box. In order to use either of these routines, the user parameter `KEEP_CUTS` must be turned on. Both point assign and box assign now work for points or boxes anywhere in space even if they are exterior to the original bounding box. If a part is empty (due to the part being assigned zero work), it is not included in the list of parts returned by box assign. Note: the original box assign algorithm was not rigorous and may have missed parts. This version is both rigorous and fast.

The Zoltan implementation of HSFC has one parameter that can be modified by the [Zoltan_Set_Param](#) function.

This partitioning algorithm is loosely based on the 2D & 3D Hilbert tables used in the Octree partitioner and on the BSFC partitioning implementation by Andrew C. Bauer, Department of Engineering, State University of New York at Buffalo, as his summer project at SNL in 2001. The box assign algorithm is loosely based on the papers by Lawder referenced both in the developers guide and the code itself. NOTE: This code can be trivially extended to any space filling curve by providing the tables implementing the curve's state transition diagram. The only dependance on the curve is through the tables and the box assign algorithm will work for all space filling curves (if we have their tables.)

Please refer to the Zoltan Developers Guide, [Appendix: Hilbert Space Filling Curve \(HSFC\)](#) for more detailed information about these algorithms.

Method String:	HSFC
Parameters:	
<i>KEEP_CUTS</i>	Information about cuts and bounding box is necessary if the application wants to add more objects to the decomposition via calls to Zoltan_LB_Point_PP_Assign or to Zoltan_LB_Box_PP_Assign . 0 = don't keep cuts; 1 = keep cuts.
<i>REDUCE_DIMENSIONS</i>	When a 3 dimensional geometry is almost flat, it may make more sense to treat it as a 2 dimensional geometry when applying the HSFC algorithm. (Coordinate values in the omitted direction are ignored for the purposes of partitioning.) If this parameter is set to 1 , a 3 dimensional geometry will be treated as 2 dimensional if is very flat, or 1 dimensional if it very thin. And a 2 dimensional geometry will be treated as 1 dimensional if it is very thin. Turning this parameter on removes the possibility that disconnected parts will appear on the surface of a flat 3 dimensional object. If the REDUCE_DIMENSIONS parameter is set, then this parameter determines when a

DEGENERATE_RATIO geometry is considered to be flat. A bounding box which is oriented to the geometry is constructed, and the lengths of its sides are tested against a ratio of 1 : **DEGENERATE_RATIO**.

Default:

KEEP_CUTS = 0
REDUCE_DIMENSIONS = 0
DEGENERATE_RATIO = 10

Required Query Functions:

[ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)
[ZOLTAN_NUM_GEOM_FN](#)
[ZOLTAN_GEOM_MULTI_FN](#) or [ZOLTAN_GEOM_FN](#)

Refinement Tree Partitioning (REFTREE)

The refinement tree based partitioning method is due to William Mitchell of the National Institute of Standards and Technology [[Mitchell](#)]. It is closely related to the Octree and Space-Filling Curve methods, except it uses the tree that represents the adaptive refinement process that created the grid. This tree is constructed through the tree-based query functions.

Each node of the refinement tree corresponds to an element that occurred during the grid refinement process. The first level of the tree (the children of the root of the tree) corresponds to the initial coarse grid, one tree node per initial element. It is assumed that the initial coarse grid does not change through the execution of the program, except that the local IDs, assignment of elements to processors, and weights can change. If any other aspect of the coarse grid changes, then the Zoltan structure should be destroyed and recreated. The children of a node in the tree correspond to the elements that were created when the corresponding element was refined. The children are ordered such that a traversal of the tree creates a space-filling curve within each initial element. If the initial elements can be ordered with a contiguous path through them, then the traversal creates a space-filling curve through all the elements. Each element has a designated "in" vertex and "out" vertex, with the out vertex of one element being the same as the in vertex of the next element in the path, in other words the path goes through a vertex to move from one element to the next (and does not go out the same vertex it came in).

The user may allow Zoltan to determine the order of the coarse grid elements, or may specify the order, which might be faster or produce a better path. If Zoltan determines the order, the user can select between an order that will produce connected parts, an order based on a Hilbert Space Filling Curve, or an order based on a Sierpinski Space Filling Curve. See the parameter REFTREE_INITPATH below. If the user provides the order, then the in/out vertices must also be supplied. Similarly, the user may specify the order and in/out vertices of the child elements, or allow Zoltan to determine them. If the user knows how to provide a good ordering for the children, this may be significantly faster than the default general algorithm. However, accelerated forms of the ordering algorithm are provided for certain types of refinement schemes and should be used in those cases. See [ZOLTAN_CHILD_LIST_FN](#). If the user always specifies the order, then the vertices and in/out vertices are not used and do not have to be provided.

Weights are assigned to the nodes of the tree. These weights need not be only on the leaves (the elements of the final grid), but can also be on interior nodes (for example, to represent work on coarse grids of a multigrid algorithm). The default weights are 1.0 at the leaves and 0.0 at the interior nodes, which produces a partition based on the number of elements in each part. An initial tree traversal is used to sum the weights, and a second traversal to cut the space-filling curve into appropriately-sized pieces and assign elements to parts. The number of parts is not necessarily equal to the number of processors.

The following limitations should be removed in the future.

- For multicomponent weights, only the first component is used.
- Heterogeneous architectures are not supported, in the sense that the computational load is equally divided over the processors. A vector of relative part sizes is used to determine the weight assigned to each part, but they are currently all equal. In the future they should be input to reflect heterogeneity.

Method String: REFTREE

Parameters:

The size of the hash table to map from global IDs to refinement tree nodes. Larger values REFTREE_HASH_SIZE require more memory but may reduce search time.

Default:

REFTREE_HASH_SIZE = 16384

Determines the method for finding an order of the elements in the initial grid.

REFTREE_INITPATH "SIERPINSKI" uses a Sierpinski Space Filling Curve and is most appropriate for grids consisting of triangles. It is currently limited to 2D.
"HILBERT" uses a Hilbert Space Filling Curve and is most appropriate for grids consisting of quadralaterals or hexahedra.
"CONNECTED" attempts to produce connected parts (guaranteed for triangles and tetrahedra), however they tend to be stringy, i.e., less compact than the SFC methods. It is most appropriate when connected parts are required.
An invalid character string will invoke the default method.

Default:

REFTREE_INITPATH = "SIERPINSKI" if the grid contains only triangles
REFTREE_INITPATH = "HILBERT" otherwise

NOTE: In Zoltan versions 1.53 and earlier the default was "CONNECTED". To reproduce old results, use *REFTREE_INITPATH* = "CONNECTED".

Required Query Functions:

[ZOLTAN_NUM_COARSE_OBJ_FN](#)
[ZOLTAN_COARSE_OBJ_LIST_FN](#) or
[ZOLTAN_FIRST_COARSE_OBJ_FN/ZOLTAN_NEXT_COARSE_OBJ_FN](#) pair
[ZOLTAN_NUM_CHILD_FN](#)
[ZOLTAN_CHILD_LIST_FN](#)
[ZOLTAN_CHILD_WEIGHT_FN](#)

The following functions are needed only if the order of the initial elements will be determined by a space filling curve method:

[ZOLTAN_NUM_GEOM_FN](#)
[ZOLTAN_GEOM_MULTI_FN](#) or [ZOLTAN_GEOM_FN](#)

Hypergraph partitioning

Hypergraph partitioning is a useful partitioning and load balancing method when connectivity data is available. It can be viewed as a more sophisticated alternative to the traditional graph partitioning.

A hypergraph consists of vertices and hyperedges. A hyperedge connects one or more vertices. A graph can be cast as a hypergraph in one of two ways: either every pair of neighboring vertices form a hyperedge, or a vertex and all its neighbors form a hyperedge. The hypergraph model is well suited to parallel computing, where vertices correspond to data objects and hyperedges represent the communication requirements. The basic partitioning problem is to partition the vertices into k approximately equal sets such that the number of cut hyperedges is minimized. Most partitioners (including Zoltan-PHG) allows a more general model where both vertices and hyperedges can be assigned weights. It has been shown that the hypergraph model gives a more accurate representation of communication cost (volume) than the graph model. In particular, for sparse matrix-vector multiplication, the hypergraph model **exactly** represents communication volume. Sparse matrices can be partitioned either along rows or columns; in the row-net model the columns are vertices and each row corresponds to an hyperedge, while in the column-net model the roles of vertices and hyperedges are reversed.

Zoltan contains a native parallel hypergraph partitioner, called PHG (Parallel HyperGraph partitioner). In addition, Zoltan provides access to [PaToH](#), a serial hypergraph partitioner. Note that PaToH is not part of Zoltan and should be obtained separately from the [PaToH web site](#). Zoltan-PHG is a fully parallel multilevel hypergraph partitioner. For further technical description, see [\[Devine et al, 2006\]](#).

A new feature available in Zoltan 3.0 is the ability to assign selected objects (vertices) to a particular part ("fixed vertices"). When objects are fixed, Zoltan will not migrate them out of the user assigned part. See the descriptions of the [ZOLTAN_NUM_FIXED_OBJ_FN](#) and [ZOLTAN_FIXED_OBJ_LIST_FN](#) query functions for a discussion of how you can define these two functions to fix objects to parts. Both PHG and PaToH support this feature.

For applications that already use Zoltan to do graph partitioning, it is easy to upgrade to hypergraph partitioning. For many applications, the hypergraph model is superior to the graph model, but in some cases the graph model should be preferred. PHG can also be used as a pure graph partitioner. See the section [graph vs. hypergraph](#) partitioning for further details.

Method String:	HYPERGRAPH
Parameters:	
<i>HYPERGRAPH_PACKAGE</i>	The software package to use in partitioning the hypergraph. <i>PHG (Zoltan, the default)</i> <i>PATOH</i>

PHG - Parallel Hypergraph and Graph Partitioning with Zoltan

This is the built-in parallel hypergraph partitioner in Zoltan.

Value of LB_METHOD: **HYPERGRAPH** (for [hypergraph](#) partitioning) or **GRAPH** (for [graph](#) partitioning)

Value of HYPERGRAPH_PACKAGE: **PHG**

Parameters:

<i>LB_APPROACH</i>	The load balancing approach: <i>REPARTITION</i> - partition but try to stay close to the current partition/distribution <i>REFINE</i> - refine the current partition/distribution; assumes only small changes <i>PARTITION</i> - partition from scratch, not taking the current data distribution into account
<i>PHG_REPART_MULTIPLIER</i>	For repartitioning, this parameter determines the trade-off between application communication (as represented by cut edges) and data migration related to rebalancing. PHG attempts to minimize the function (PHG_REPART_MULTIPLIER* edge_cut + migration volume). The migration volume is measured using the ZOLTAN_OBJ_SIZE_MULTI_FN or ZOLTAN_OBJ_SIZE_FN query functions. Make sure the units for edges and object sizes are the same. Simply put, to emphasize communication within the application, use a large value for PHG_REPART_MULTIPLIER. Typically this should be proportional to the number of iterations between load-balancing calls.
<i>PHG_MULTILEVEL</i>	This parameter specifies whether a multilevel method should be used (1) or not (0). Multilevel methods produce higher quality but require more execution time and memory.
<i>PHG_EDGE_WEIGHT_OPERATION</i>	Operation to be applied to edge weights supplied by different processes for the same hyperedge: <i>ADD</i> - the hyperedge weight will be the sum of the supplied weights <i>MAX</i> - the hyperedge weight will be the maximum of the supplied weights <i>ERROR</i> - if the hyperedge weights are not equal, Zoltan will flag an error, otherwise the hyperedge weight will be the value returned by the processes
<i>ADD_OBJ_WEIGHT</i>	Add another object (vertex) weight. Currently multi-weight partitioning is not supported, but this parameter may also be used for implicit vertex weights. Valid values are: <i>NONE</i> <i>UNIT</i> or <i>VERTICES</i> (each vertex has weight 1.0) <i>PINS</i> or <i>NONZEROS</i> or <i>vertex degree</i> (vertex weight equals number of hyperedges containing it, i.e., the degree)
<i>PHG_CUT_OBJECTIVE</i>	Selects the partitioning objective, <i>CONNECTIVITY</i> or <i>HYPEREDGES</i> . While hyperedges simply counts the number of hyperedges cut, the connectivity metric weights each cut edge by the number of participating processors - 1 (aka the k-1 cut metric). The connectivity metric better represents communication volume for most applications. The hyperedge metric is useful for certain applications, e.g., minimizing matrix border size in block matrix decompositions.

<i>PHG_OUTPUT_LEVEL</i>	Level of verbosity; 0 is silent.
<i>CHECK_HYPERGRAPH</i>	Check that the query functions return valid input data; 0 or 1. (This slows performance; intended for debugging.)
<i>PHG_COARSENING_METHOD</i>	Low-level parameter: The method to use in the matching/coarsening phase: <i>AGG</i> - agglomerative inner product matching (a.k.a. agglomerative heavy connectivity matching); gives high quality. <i>IPM</i> - inner product matching (a.k.a. heavy connectivity matching); gives high quality. <i>L-IPM</i> - local IPM on each processor. Faster but usually gives poorer quality. <i>A-IPM</i> - alternate between IPM and L-IPM. (A compromise between speed and quality.) <i>none</i> - no coarsening
<i>PHG_COARSEPARTITION_METHOD</i>	Low-level parameter: Method to partition the coarsest (smallest) hypergraph; typically done in serial: <i>RANDOM</i> - random <i>LINEAR</i> - linear assignment of the vertices (ordered by the user query function) <i>GREEDY</i> - greedy method based on minimizing cuts <i>AUTO</i> - automatically select from the above methods (in parallel, the processes will do different methods)
<i>PHG_REFINEMENT_METHOD</i>	Low-level parameter: Refinement algorithm: <i>FM</i> - approximate Fiduccia-Mattheyses (FM) <i>NO</i> - no refinement
<i>PHG_REFINEMENT_QUALITY</i>	Low-level parameter: Knob to control the trade-off between run time and quality. 1 is the recommended (default) setting, >1 gives more refinement (higher quality partitions but longer run time), while <1 gives less refinement (and poorer quality).
<i>PHG_RANDOMIZE_INPUT</i>	Low-level parameter: Randomize layout of vertices and hyperedges in internal parallel 2D layout? Setting this parameter to 1 often reduces Zoltan-PHG execution time.
<i>PHG_EDGE_SIZE_THRESHOLD</i>	Low-level parameter: Value 0.0 through 1.0, if number of vertices in hyperedge divided by total vertices in hypergraph exceeds this fraction, the hyperedge will be omitted.
<i>PHG_PROCESSOR_REDUCTION_LIMIT</i>	Low-level parameter: In V-cycle, redistribute coarsened hypergraph to this fraction of processors when number of pins in coarsened hypergraph is less than this fraction of original number of pins. Original number of pins is redefined after each processor redistribution.

Default values:

LB_APPROACH = *REPARTITION*
PHG_REPART_MULTIPLIER=100
PHG_MULTILEVEL=1 if *LB_APPROACH* = *partition* or *repartition*; 0 otherwise.
PHG_EDGE_WEIGHT_OPERATION=max
ADD_OBJ_WEIGHT=none
PHG_CUT_OBJECTIVE=connectivity
CHECK_HYPERGRAPH=0
PHG_OUTPUT_LEVEL=0

PHG_COARSENING_METHOD=agg
PHG_COARSEPARTITION_METHOD=auto
PHG_REFINEMENT_METHOD=fm
PHG_REFINEMENT_QUALITY=1
PHG_RANDOMIZE_INPUT=0
PHG_EDGE_SIZE_THRESHOLD=0.25
PHG_PROCESSOR_REDUCTION_LIMIT=0.5

**Required Query Functions for
*LB_METHOD = HYPERGRAPH:***

[ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)
[ZOLTAN_HG_SIZE_CS_FN](#)
[ZOLTAN_HG_CS_FN](#)

**Optional Query Functions for
*LB_METHOD = HYPERGRAPH:***

[ZOLTAN_OBJ_SIZE_MULTI_FN](#) or [ZOLTAN_OBJ_SIZE_FN](#) for
LB_APPROACH=Repartition.
[ZOLTAN_PART_MULTI_FN](#) or [ZOLTAN_PART_FN](#) for
LB_APPROACH=Repartition and for [REMAP=1](#).
[ZOLTAN_HG_SIZE_EDGE_WTS_FN](#)
[ZOLTAN_HG_EDGE_WTS_FN](#)
[ZOLTAN_NUM_FIXED_OBJ_FN](#)
[ZOLTAN_FIXED_OBJ_LIST_FN](#)

**Note for
*LB_METHOD = HYPERGRAPH:***

It is possible to provide the graph query functions instead of the hypergraph queries, though this is not recommended. If only graph query functions are registered, Zoltan will automatically create a hypergraph from the graph, but this is not equivalent to graph partitioning. In particular, the edge weights will not be accurate.

**Required Query Functions for
*LB_METHOD = GRAPH:***

[ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)
[ZOLTAN_NUM_EDGES_MULTI_FN](#) or
[ZOLTAN_NUM_EDGES_FN](#)
[ZOLTAN_EDGE_LIST_MULTI_FN](#) or [ZOLTAN_EDGE_LIST_FN](#)

**Optional Query Functions for
*LB_METHOD = GRAPH:***

[ZOLTAN_OBJ_SIZE_MULTI_FN](#) or [ZOLTAN_OBJ_SIZE_FN](#) for
LB_APPROACH=Repartition.
[ZOLTAN_PART_MULTI_FN](#) or [ZOLTAN_PART_FN](#) for
LB_APPROACH=Repartition and for [REMAP=1](#).

PaToH

[PaToH](#) is a serial hypergraph partitioning package. It is not distributed with Zoltan and must be obtained separately. (You will also need to modify your Zoltan configuration file after installation.)

Since PaToH is serial, it can only be used in Zoltan on a single processor. PaToH is faster than Zoltan PHG in serial mode, and often produces (slightly) better partition quality.

Value of LB_METHOD: **HYPERGRAPH**

**Value of
HYPERGRAPH_PACKAGE:** **PATOH**

Parameters:

<i>PATOH_ALLOC_POOL0</i>	<i>non-zero</i> (the value of the MemMul_CellNet PATOH parameter) <i>0</i> (the default)
<i>PATOH_ALLOC_POOL1</i>	<i>non-zero</i> (the value of the MemMul_Pins PATOH parameter) <i>0</i> (the default)
<i>USE_TIMERS</i>	<i>1</i> (time operations and print results) <i>0</i> (don't time, the default)
<i>PHG_EDGE_SIZE_THRESHOLD</i>	Value 0.0 through 1.0, if number of vertices in hyperedge divided by total vertices in hypergraph exceeds this fraction, the hyperedge will be omitted.
<i>ADD_OBJ_WEIGHT</i>	Add implicit vertex (object) weight. Multi-weight partitioning is not yet supported, so currently either the user-defined weight or the implicit weight will be used. Valid values: <i>none</i> (the default if OBJ_WEIGHT_DIM > 0) <i>unit</i> or <i>vertices</i> (each vertex has weight 1.0, default if OBJ_WEIGHT_DIM is zero) <i>pins</i> or <i>nonzeros</i> or <i>vertex degree</i> (vertex weight equals number hyperedges containing it)
<i>CHECK_HYPERGRAPH</i>	Check that the query functions return valid input data; 0 or 1. (This slows performance; intended for debugging.)

Default Values:

PATOH_ALLOC_POOL0 = 0
PATOH_ALLOC_POOL1 = 0
USE_TIMERS = 0
PHG_EDGE_SIZE_THRESHOLD = 0.25
ADD_OBJ_WEIGHT = none
CHECK_HYPERGRAPH = 0

Required Query Functions:

[ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)
[ZOLTAN_HG_SIZE_CS_FN](#)
[ZOLTAN_HG_CS_FN](#)

Optional Query Functions:

[ZOLTAN_HG_SIZE_EDGE_WTS_FN](#)
[ZOLTAN_HG_EDGE_WTS_FN](#)

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Note: See also [hypergraph partitioning](#).

Graph partitioning

Zoltan performs graph partitioning when the *LB_METHOD* parameter is set to GRAPH. Zoltan provides three packages capable of partitioning a graph. The package is chosen by setting the GRAPH_PACKAGE parameter. Two packages (ParMetis and Scotch) are external packages and not part of Zoltan but accessible via Zoltan. The last package is PHG, Zoltan's native hypergraph partitioner. PHG will treat the graph as a regular hypergraph with edge size two. Since PHG was designed for general hypergraphs, it is usually slower than graph partitioners but often produces better quality.

Method String:	GRAPH
Parameters:	
<i>GRAPH_PACKAGE</i>	The software package to use in partitioning the graph. <i>PHG</i> (<i>default</i>) <i>ParMETIS</i> <i>Scotch/PT-Scotch</i>

Graph vs Hypergraph Partitioning

Graph partitioning has proven quite useful in scientific computing. [Hypergraph partitioning](#) is a more recent improvement that uses a hypergraph model, which is often a more accurate model than the graph model for scientific computing. (Hypergraphs contain hyperedges which connect two *or more* vertices.) See [\[Catalyurek & Aykanat\]](#) and [\[Hendrickson & Kolda\]](#) for further details. You do not need to understand the underlying models to use graph or hypergraph partitioning for load-balancing in Zoltan. The basic trade-offs are:

- Hypergraph partitioning usually produces partitions (assignments) of higher quality than graph partitioning, which may reduce communication time in parallel applications (up to 30-40% reduction has been reported). However, hypergraph partitioning takes longer time to compute.
- The graph model is restricted to symmetric data dependencies. If you have a non-symmetric problem, we recommend hypergraph partitioning.

Migrating from ParMetis to PHG in Zoltan

If you already use Zoltan for graph partitioning (via ParMetis), there are three ways to switch to the Zoltan-PHG hypergraph partitioner:

1. The quick and easy way: Just change the [LB_METHOD](#) to "Hypergraph". Zoltan will then use the graph query functions (presumably already implemented) to construct a hypergraph model, which is similar to but not equivalent to the graph.
2. The proper way: Change the LB_METHOD, but also implement and register the [hypergraph query functions](#) required by Zoltan. These may give a more accurate representation of data dependencies (and communication requirements) for your application.
3. If you really want graph (not hypergraph) partitioning: Just change the [LB_METHOD](#) to "Graph". Zoltan will then use PHG as a graph partitioner, which is slower than ParMetis but often produces better partitions (lower cuts).

Technical note: A hypergraph is constructed from the graph as follows: The vertices are the same in the hypergraph as in the graph. For each vertex v , create a hyperedge that consists of all neighbors in the graph and v itself.

ParMETIS - Parallel Graph Partitioning

[ParMETIS](#) is a parallel library for graph partitioning (for static load balancing) and repartitioning (for dynamic load balancing) developed at the University of Minnesota by Karypis, Schloegel and Kumar [[Karypis and Kumar](#)]. ParMETIS is therefore strictly speaking not a method but rather a collection of methods. In the Zoltan context, ParMETIS is a method with many sub-methods. Zoltan provides an interface to all the ParMETIS (sub-)methods. The user selects which ParMETIS method to use through the parameter `PARMETIS_METHOD`. Most of the ParMETIS methods are based on either multilevel Kernighan-Lin partitioning or a diffusion algorithm. The names of the ParMETIS methods used by Zoltan are identical to those in the ParMETIS library. For further information about the various [ParMETIS](#) methods and parameters, please consult the [ParMETIS](#) User's Guide.

Graph partitioning is a useful abstraction for load balancing. The main idea is to represent the computational application as a weighted graph. The nodes or vertices in the graph correspond to objects in Zoltan. Each object may have a weight that normally represents the amount of computation. The edges or arcs in the graph usually correspond to communication costs. In graph partitioning, the problem is to find a partition of the graph (that is, each vertex is assigned to one out of k possible sets called parts) that minimizes the cut size (weight) subject to the parts having approximately equal size (weight). In repartitioning, it is assumed that a partition already exists. The problem is to find a good partition that is also "similar" in some sense to the existing partition. This keeps the migration cost low. All the problems described above are NP-hard so no efficient exact algorithm is known, but heuristics work well in practice.

We give only a brief summary of the various ParMETIS methods here; for more details see the [ParMETIS](#) documentation. The methods fall into three categories:

1. Part* - Perform graph partitioning without consideration of the initial distribution. (`LB_APPROACH=partition`)
2. AdaptiveRepart - Incremental algorithms with small migration cost. (`LB_APPROACH=repartition`)
3. Refine* - Refines a given partition (balance). Can be applied multiple times to reduce the communication cost (cut weight) if desired. (`LB_APPROACH=refine`)

As a rule of thumb, use one of the Part* methods if you have a poor initial balance and you are willing to spend some time doing migration. One such case is static load balancing; that is, you need to balance only once. Use AdaptiveRepart or the Repart* methods when you have a reasonably good load balance that you wish to update incrementally. These methods are well suited for dynamic load balancing (for example, adaptive mesh refinement). A reasonable strategy is to call PartKway once to obtain a good initial balance and later update this balance using AdaptiveRepart.

Zoltan is currently compatible with ParMETIS version 3.1. (or 3.1.1) (Version 2.0 is obsolete and is no longer supported by Zoltan.) The ParMETIS source code can be obtained from the [ParMETIS home page](#). As a courtesy service, a recent, compatible version of the ParMETIS source code is distributed with Zoltan. However, ParMETIS is a completely separate library. If you do not wish to install ParMETIS, it is possible to compile Zoltan without any references to ParMETIS (when you 'make' Zoltan, comment out the `PARMETIS_LIBPATH` variable in the configuration file [Utilities/Config/Config.<platform>](#)).

Zoltan supports the multiconstraint partitioning feature of ParMETIS through multiple object weights (see [OBJ_WEIGHT_DIM](#)).

The graph given to Zoltan/ParMETIS must be symmetric. Any self edges (loops) will be ignored. Multiple edges between a pair of vertices is not allowed. All weights must be non-negative. The graph does not have to be connected.

Value of `LB_METHOD`: **GRAPH**
Value of
`GRAPH_PACKAGE`: **Parmetis**

Parameters:

<i>LB_APPROACH</i>	The load balancing approach: <i>PARTITION</i> - partition from scratch, not taking the current data distribution into account <i>REPARTITION</i> - partition but try to stay close to the current partition/distribution <i>REFINE</i> - refine the current partition/distribution; assumes only small changes
<i>PARMETIS_METHOD</i>	The specific ParMETIS method to be used (see below). Note: See also LB_APPROACH , which is a simpler way to specify the overall load balance approach. Only use <i>PARMETIS_METHOD</i> if you really need a specific implementation. <i>PartKway</i> - multilevel Kernighan-Lin partitioning <i>PartGeom</i> - space filling curves (coordinate based) <i>PartGeomKway</i> - hybrid method based on PartKway and PartGeom (needs both graph data and coordinates) <i>AdaptiveRepart</i> - adaptive repartitioning (only in ParMETIS 3.0 and higher) <i>RefineKway</i> - refine the current partition (balance) The method names are case insensitive.
<i>PARMETIS_OUTPUT_LEVEL</i>	Amount of output the load-balancing algorithm should produce. 0 = no output, 1 = print timing info. Turning on more bits displays more information (for example, 3=1+2, 5=1+4, 7=1+2+4).
<i>PARMETIS_COARSE_ALG</i>	Coarse algorithm for PartKway. 1 = serial, 2 = parallel. (ParMETIS 2 only)
<i>PARMETIS_SEED</i>	Random seed for ParMETIS.
<i>PARMETIS_ITR</i>	Ratio of interprocessor communication time to redistribution time. A high value will emphasize reducing the edge cut, while a small value will try to keep the change in the new partition (distribution) small. This parameter is used only by AdaptiveRepart. A value of between 100 and 1000 is good for most problems.
<i>CHECK_GRAPH</i>	Level of error checking for graph input: 0 = no checking, 1 = on-processor checking, 2 = full checking. (CHECK_GRAPH==2 is very slow and should be used only during debugging).
<i>SCATTER_GRAPH</i>	Scatter graph data by distributing contiguous chunks of objects (vertices) of roughly equal size to each processor before calling the partitioner. 0 = don't scatter; 1 = scatter only if all objects are on a single processor; 2 = scatter if at least one processor owns no objects; 3 = always scatter.
<i>GRAPH_SYMMETRIZE</i>	How to symmetrize the graph: NONE = graph is symmetric and no symmetrization is needed TRANSPOSE = if M is adjacency matrix of the input graph, output will be the graph representation of $M+M^T$ BIPARTITE = graph is symmetrized in a bipartite way : $\begin{bmatrix} 0 & M \\ M^t & 0 \end{bmatrix}$
<i>GRAPH_SYM_WEIGHT</i>	How edge weights are handled during symmetrization: ADD = weights of each arc are added MAX = only the heaviest arc weight is kept

See more informations about graph build options on this [page](#)

Default values:

LB_APPROACH = Repartition
PARMETIS_METHOD = AdaptiveRepart
PARMETIS_OUTPUT_LEVEL = 0
PARMETIS_COARSE_ALG = 2
PARMETIS_SEED = 15

PARMETIS_ITR = 100
USE_OBJ_SIZE = 1
CHECK_GRAPH = 1
SCATTER_GRAPH = 1
GRAPH_SYMMETRIZE = NONE
GRAPH_SYM_WEIGHT = ADD

Required Query Functions:

For all submethods: [ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)

Only PartGeom &
PartGeomKway: [ZOLTAN_NUM_GEOM_FN](#)

[ZOLTAN_GEOM_MULTI_FN](#) or [ZOLTAN_GEOM_FN](#)
All but PartGeom: [ZOLTAN_NUM_EDGES_MULTI_FN](#) or [ZOLTAN_NUM_EDGES_FN](#)
[ZOLTAN_EDGE_LIST_MULTI_FN](#) or [ZOLTAN_EDGE_LIST_FN](#)

Optional Query Functions:

[ZOLTAN_OBJ_SIZE_MULTI_FN](#) or [ZOLTAN_OBJ_SIZE_FN](#) for
PARMETIS_METHOD=*AdaptiveRepart*
[ZOLTAN_PART_MULTI_FN](#) or [ZOLTAN_PART_FN](#) for part remapping in
ParMETIS.

Graph partitioning: Scotch/PT-Scotch

For more information about graph partitioning, see [ParMetis](#).

Scotch is a library for ordering and partitioning, developed at LaBRI in Bordeaux, France. PT-Scotch is the parallel module in Scotch. (We use the name PT-Scotch when it applies only to parallel version, Scotch when it concerns at least the serial version.) Scotch is a third-party library in Zoltan and should be obtained separately from the [Scotch web site](#). Currently, Zoltan supports version 5.1 of Scotch, compiled with the support of PT-Scotch and *int* coding for vertices, not *long* (on architectures where *sizeof(int) != sizeof(long)*).

If the parameter [SCOTCH_TYPE](#) is set to LOCAL and the number of processors is 1 (e.g., `mpirun -np 1`), the sequential version of Scotch is called.

Both the serial (Scotch) and the parallel (PT-Scotch) compute k-way partitioning by doing recursive bisection.

Scotch must be used in the context `LB_APPROACH=partition`, to balance poorly distributed data. Unlike ParMetis, Scotch doesn't support the multiconstraint partitioning feature. However, for single constraint partitioning it seems to produce better results than ParMetis although it requires less memory.

Value of LB_METHOD:	GRAPH
Value of GRAPH_PACKAGE:	Scotch
Parameters:	
<i>LB_APPROACH</i>	The load balancing approach: <i>PARTITION</i> - partition from scratch, not taking the current data distribution into account
<i>SCOTCH_METHOD</i>	For now, always set to RBISECT
<i>SCOTCH_STRAT</i>	The Scotch strategy string; see the Scotch documentation.
<i>SCOTCH_STRAT_FILE</i>	A file that contains an arbitrary long Scotch strategy string; see the Scotch documentation.
<i>SCOTCH_TYPE</i>	Whether or not the parallel library is used. The default value is <i>GLOBAL</i> and if not set to <i>LOCAL</i> , the parallel graph partitioning is used.
<i>CHECK_GRAPH</i>	Level of error checking for graph input: 0 = no checking, 1 = on-processor checking, 2 = full checking. (<code>CHECK_GRAPH==2</code> is very slow and should be used only during debugging).
<i>SCATTER_GRAPH</i>	Scatter graph data by distributing contiguous chunks of objects (vertices) of roughly equal size to each processor before calling the partitioner. 0 = don't scatter; 1 = scatter only if all objects are on a single processor; 2 = scatter if at least one processor owns no objects; 3 = always scatter.
<i>GRAPH_SYMMETRIZE</i>	How to symmetrize the graph: NONE = graph is symmetric and no symmetrization is needed TRANSPOSE = if M is adjacency matrix of the input graph, output will be the graph representation of $M+M^T$ BIPARTITE = graph is symmetrized in a bipartite way : $\begin{bmatrix} 0 & M \\ M^t & 0 \end{bmatrix}$
<i>GRAPH_SYM_WEIGHT</i>	How edge weights are handled during symmetrization: ADD = weights of each arc are added MAX = only the heaviest arc weight is kept

See more informations about graph build options on this [page](#)

Default values:

LB_APPROACH = Partition
SCOTCH_METHOD = RBISECT
SCOTCH_STRAT = "" (default Scotch strategy)
SCOTCH_STRAT_FILE = "" (default Scotch strategy)
CHECK_GRAPH = 1
SCATTER_GRAPH = 1
GRAPH_SYMMETRIZE = NONE
GRAPH_SYM_WEIGHT = ADD

**Required Query
Functions:**

[ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)
[ZOLTAN_NUM_EDGES_MULTI_FN](#) or [ZOLTAN_NUM_EDGES_FN](#)
[ZOLTAN_EDGE_LIST_MULTI_FN](#) or [ZOLTAN_EDGE_LIST_FN](#)

Hierarchical Partitioning (HIER)

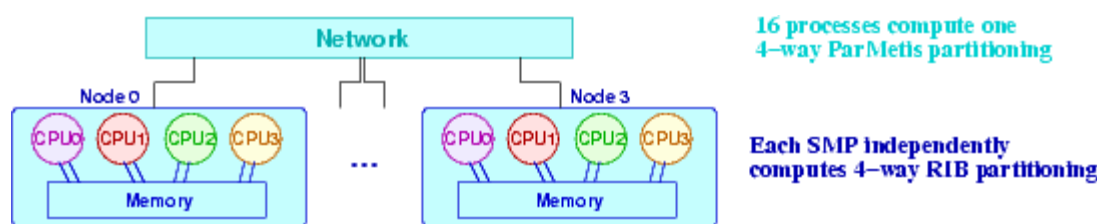
Hierarchical partitioning allows the creation of hybrid partitions, computed using combinations of other Zoltan procedures. For hierarchical and heterogeneous systems, different choices may be appropriate in different parts of the parallel environment. There are tradeoffs in execution time and partition quality (*e.g.*, surface indices/communication volume, interprocess connectivity, strictness of load balance) and some may be more important than others in some circumstances. For example, consider a cluster of symmetric multiprocessor (SMP) nodes connected by Ethernet. A more costly graph partitioning can be done to partition among the nodes, to minimize communication across the slow network interface, possibly at the expense of some computational imbalance. Then, a fast geometric algorithm can be used to partition independently within each node.

Zoltan's hierarchical balancing, implemented by Jim Teresco (Williams College) during a 2003-04 visit to Sandia, automates the creation of hierarchical partitions [[Teresco, et al.](#)]. It can be used directly by an application or be guided by the tree representation of the computational environment created and maintained by the [Dynamic Resource Utilization Model \(DRUM\)](#) [[Devine, et al.](#), [Faik, et al.](#), [Teresco, et al.](#)].

The hierarchical balancing implementation utilizes a lightweight intermediate structure and a set of callback functions that permit an automated and efficient hierarchical balancing which can use any of the procedures available within Zoltan without modification and in any combination. Hierarchical balancing is invoked by an application the same way as other Zoltan procedures and interfaces with applications through callback functions. A hierarchical balancing step begins by building an intermediate structure using these callbacks. This structure is an augmented version of the graph structure that Zoltan builds to make use of the ParMetis and Jostle libraries. The hierarchical balancing procedure then provides its own callback functions to allow existing Zoltan procedures to be used to query and update the intermediate structure at each level of a hierarchical balancing. After all levels of the hierarchical balancing have been completed, Zoltan's usual migration arrays are constructed and returned to the application. Thus, only lightweight objects are migrated internally between levels, not the (larger and more costly) application data.

Hierarchical partitioning requires three callback functions to specify the number of levels ([ZOLTAN_HIER_NUM_LEVELS_FN](#)), which parts each process should compute at each level ([ZOLTAN_HIER_PART_FN](#)), and which method and parameters to be used at each level ([ZOLTAN_HIER_METHOD_FN](#)). These are in addition to the callbacks needed to specify objects, coordinates, and graph edges. This fairly cumbersome interface can be avoided by using the separately available [zoltanParams](#) library. This allows a file-based description to replace these callbacks. A more direct interface with DRUM's hierarchical machine model is also planned, allowing hierarchical balancing parameters to be set by a graphical configuration tool.

We use a simple example to illustrate the use of the callback mechanism to specify hierarchical a hierarchical partitioning. In the figure [below](#), a hierarchical computing environment and a desired hierarchical partitioning is shown.



Assume we start one process for each processor, with the processes of ranks 0-3 assigned to Node 0, 4-7 to Node 1, 8-11 to Node 2, and 12-15 to Node 3. When hierarchical partitioning is invoked, the following callbacks will be made, and the following actions should be taken by the callbacks on each node.

1. The [ZOLTAN_HIER_NUM_LEVELS_FN](#) callback is called. All processes should return 2, the number of

levels in the hierarchy.

2. The **ZOLTAN_HIER_PART_FN** callback is called, with a **level** parameter equal to 0. This means the callback should return, on each process, the part number to be computed at level 0 by that process. Since in our example, the 16 processes are computing a four-way partition at level 0, processes with ranks 0-3 should return 0, ranks 4-7 should return 1, 8-11 should return 2, and 12-15 should return 3.
3. The **ZOLTAN_HIER_METHOD_FN** callback is called, with a **level** parameter equal to 0, and the Zoltan_Struct that has been allocated internally by the hierarchical partitioning procedure is passed as **zz**. The callback should use Zoltan_Set_Param to specify an LB_METHOD and any other parameters that should be done by the four-way partition across the 16 processes at level 0. In this case, two calls might be appropriate:

```
Zoltan_Set_Param(zz, "LB_METHOD", "PARMETIS");  
Zoltan_Set_Param(zz, "PARMETIS_METHOD", "PARTKWAY");
```

At this point, Zoltan's hierarchical balancing procedure can proceed with the level 0 partition, using ParMetis' PARTKWAY method to produce a four-way partition across the 16 processes, with part 0 distributed among the processes with ranks 0-3, part 1 distributed among 4-7, part 2 distributed among 8-11, and part 3 distributed among 12-15.

4. The **ZOLTAN_HIER_PART_FN** callback is called again, this time with **level** equal to 1. At level 1, each group of four processes on the same node is responsible for computing its own four-way partition. To specify this, processes with ranks 0, 4, 8, and 12 should return 0, ranks 1, 5, 9, and 13 should return 1, ranks 2, 6, 10, and 14 should return 2, and ranks 3, 7, 11, and 15 should return 3. Note that we are specifying four separate four-way partitions at this level, so each group of four processes on the same node will have one process computing each part 0, 1, 2, and 3, for that group.
5. The **ZOLTAN_HIER_METHOD_FN** callback is called again, with **level** equal to 1, and another internally-allocated Zoltan_Struct passed as **zz**. Here, we want all processes to be computing a partition using recursive inertial bisection. The following call would be appropriate inside the callback:

```
Zoltan_Set_Param(zz, "LB_METHOD", "RIB");
```

Additional Zoltan_Set_Param calls would be used to specify any additional procedures. Note that in this case, we are computing four separate partitions but all with the same LB_METHOD. It would be allowed to specify different LB_METHODs for each group, but all processes cooperating on a partition must agree on their LB_METHOD and other parameters (just like any other Zoltan partitioning).

At this point, Zoltan's hierarchical balancing procedure can proceed with the level 1 partition, using four independent recursive inertial bisections produce the four four-way partitions across the processes on each node. Since this is the final level, the 16 resulting parts are returned by the hierarchical balancing procedure to the calling application.

Method String:

HIER

Parameters:

HIER_CHECKS

If set to 1, forces "sanity checks" to be performed on the intermediate structure when it is first created, and after the partitioning at each level.

HIER_DEBUG_LEVEL

Amount of output the hierarchical partitioning procedures should produce.
0 = no statistics; 1 = hierarchical balancing lists; 2 = all debugging information.

Default:

HIER_CHECKS = 0

HIER_DEBUG_LEVEL = 1

Required Query Functions:

[ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)
[ZOLTAN_HIER_NUM_LEVELS_FN](#),
[ZOLTAN_HIER_PART_FN](#), and
[ZOLTAN_HIER_METHOD_FN](#).
[ZOLTAN_NUM_GEOM_FN](#)
[ZOLTAN_GEOM_MULTI_FN](#) or [ZOLTAN_GEOM_FN](#)
[ZOLTAN_NUM_EDGES_MULTI_FN](#) or
[ZOLTAN_NUM_EDGES_FN](#)
[ZOLTAN_EDGE_LIST_MULTI_FN](#) or
[ZOLTAN_EDGE_LIST_FN](#)

Only if one of the methods used at some level of hierarchical partitioning requires geometric information:

Only if one of the methods used at some level of hierarchical partitioning requires graph information:

Ordering Algorithms

The following graph (sparse matrix) ordering algorithms are currently included in the Zoltan toolkit:

- [Nested dissection by METIS/ParMETIS](#)
- [Nested dissection by Scotch/PT-Scotch](#)
- [Local ordering with Hilbert space filling curves](#)

These methods produce orderings for various applications (e.g., reducing fill in sparse matrix factorizations). Ordering is accessed through calls to [Zoltan_Order](#).

Third-party libraries

Currently, most ordering in Zoltan is provided through the third-party libraries METIS/ParMETIS and PT-Scotch. The one exception is the local Hilbert space filling curve ordering. To use the other methods, a third-party library must be present.

Ordering Parameters

While the overall behavior of Zoltan is controlled by [general Zoltan parameters](#), the behavior of each ordering method is controlled by parameters specific to ordering which are also set by calls to [Zoltan_Set_Param](#). Many of these parameters are specific to individual ordering algorithms, and are listed in the descriptions of the individual algorithms. However, several have meaning across multiple ordering algorithms. These parameters are described below.

Parameters:

ORDER_METHOD The order algorithm used by Zoltan is specified by this parameter. Valid values are

- "METIS" (sequential [nodal nested dissection by METIS](#)),
- "PARMETIS" (parallel [nodal nested dissection by ParMETIS](#)),
- "SCOTCH" (sequential ordering using [Scotch](#)),
- "PTSCOTCH" (parallel ordering using [PT-Scotch](#)),
- "LOCAL_HSFC" (local ordering using [Hilbert space filling curves](#)), and
- "NONE" (for no ordering).

Default Values:

ORDER_METHOD = PARMETIS (with MPI), METIS (when no MPI)

Nested Dissection by METIS/ParMETIS

Nested Dissection (ND) is a popular method to compute fill-reducing orderings for sparse matrices. It can also be used for other ordering purposes. The algorithm recursively finds a separator (bisector) in a graph, orders the nodes in the two subsets first, and nodes in the separator last. In METIS/ParMETIS, the recursion is stopped when the graph is smaller than a certain size, and some version of minimum degree ordering is applied to the remaining graph.

METIS computes a (local) ordering of the objects (currently only for serial runs), while ParMETIS performs a global ordering of all the objects. ParMETIS currently (version 3.1, supported by Zoltan) requires that the number of processors is a power of two.

If ORDER_METHOD=PARMETIS ParMETIS is called, but if it is METIS, METIS is called.

Order_Method String: METIS or PARMETIS

Parameters:

See [ParMETIS](#). Note that the PARMETIS options are ignored when METIS is called.

Required Query Functions:

[ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)
[ZOLTAN_NUM_EDGES_MULTI_FN](#) or [ZOLTAN_NUM_EDGES_FN](#)
[ZOLTAN_EDGE_LIST_MULTI_FN](#) or [ZOLTAN_EDGE_LIST_FN](#)

Nested Dissection by Scotch

Nested Dissection is a popular method to compute fill-reducing orderings for graphs and sparse matrices. It can also be used for other ordering purposes. The algorithm recursively finds a separator (bisector) in a graph, orders the nodes in the two subsets first, and nodes in the separator last.

Scotch is a library for ordering and partitioning, developed at LaBRI in Bordeaux, France. PT-Scotch is the parallel module in Scotch. (We use the names Scotch and PT-Scotch interchangeably.) Scotch is a third-party library in Zoltan and should be obtained separately from the [Scotch web site](#). Currently, Zoltan supports version 5.1 of Scotch, compiled with the support of PT-Scotch and *int* coding for vertices, not *long* (on architectures where *sizeof(int) != sizeof(long)*).

If the parameter [ORDER_METHOD](#) is set to SCOTCH, the sequential version of Scotch is called. If the parameter [ORDER_METHOD](#) is set to PTSCOTCH, the parallel version of Scotch (PT-Scotch) is called.

Order_Method String:	PTSCOTCH or SCOTCH
Parameters:	
SCOTCH_METHOD	For now, always set to NODEND
SCOTCH_STRAT	The Scotch strategy string; see the Scotch documentation.
SCOTCH_STRAT_FILE	A file that contains an arbitrary long Scotch strategy string; see the Scotch documentation.

Required Query Functions:

[ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)
[ZOLTAN_NUM_EDGES_MULTI_FN](#) or [ZOLTAN_NUM_EDGES_FN](#)
[ZOLTAN_EDGE_LIST_MULTI_FN](#) or [ZOLTAN_EDGE_LIST_FN](#)

Coloring Algorithms

The following coloring algorithms are currently included in the Zoltan library:

[Parallel Coloring](#)

They are accessed through calls to [Zoltan_Color](#).

Coloring Parameters

While the overall behavior of Zoltan is controlled by [general Zoltan parameters](#), the behavior of each coloring method is controlled by parameters specific to coloring which are also set by calls to [Zoltan_Set_Param](#). These parameters are described below.

Parameters:

<i>DISTANCE</i>	The maximum distance between two objects that should not get the same color is specified by this parameter. Valid values are "1" (for distance-1 coloring) and "2" (for distance-2 coloring).
<i>SUPERSTEP_SIZE</i>	Number of local objects to be colored on each processor before exchanging color information. SUPERSTEP_SIZE should be greater than 0.
<i>COMM_PATTERN</i>	Valid values are "S" (synchronous) and "A" (asynchronous). If synchronous communication is selected, processors are forced to wait for the color information from all other processors to be received before proceeding with coloring of the next SUPERSTEP_SIZE number of local objects. If asynchronous communication is selected, there is no such restriction.
<i>COLOR_ORDER</i>	Valid values are "I" (internal first), "B" (boundary first) and "U" (unordered). If "I" is selected, each processor colors its internal objects before boundary objects. If "B" is selected, each processor colors its boundary objects first. If "U" is selected, there is no such distinction between internal and boundary objects. "U" is not implemented for distance-2 coloring.
<i>COLORING_METHOD</i>	Currently only "F" (first-fit) is implemented. By using "F", the smallest available color that will not cause a conflict is assigned to the object that is being colored.

Options for graph build See more informations about graph build options on this [page](#)

Default Values:

DISTANCE = 1
SUPERSTEP_SIZE = 100
COMM_PATTERN = S
COLOR_ORDER = I
COLORING_METHOD = F

Parallel Coloring

The parallel coloring algorithm in Zoltan is based on the work of [Boman et al.](#) for distance-1 coloring and [Bozdag et al.](#) for distance-2 coloring. It was implemented in Zoltan by Doruk Bozdag and Umit Catalyurek, Department of Biomedical Informatics, Ohio State University. Distance-1 coloring algorithm is an iterative data parallel algorithm that proceeds in two-phased rounds. In the first phase, processors concurrently color the vertices assigned to them. Adjacent vertices colored in the same parallel step of this phase may result in inconsistencies. In the second phase, processors concurrently check the validity of the colors assigned to their respective vertices and identify a set of vertices that needs to be re-colored in the next round to resolve the detected inconsistencies. The algorithm terminates when every vertex has been colored correctly. To reduce communication frequency, the coloring phase is further decomposed into computation and communication sub-phases. In a communication sub-phase processors exchange recent color information. During a computation sub-phase, a number of vertices determined by the SUPERSTEP_SIZE parameter, rather than a single vertex, is colored based on currently available color information. With an appropriate choice of a value for SUPERSTEP_SIZE, the number of ensuing conflicts can be kept low while at the same time preventing the runtime from being dominated by the sending of a large number of small messages. The distance-2 graph coloring problem aims at partitioning the vertex set of a graph into the fewest sets consisting of vertices pairwise at distance greater than two from each other. The algorithm is an extension of the parallel distance-1 coloring algorithm.

Parameters:

See [Coloring Algorithms](#).

Required Query Functions:

[ZOLTAN_NUM_OBJ_FN](#)
[ZOLTAN_OBJ_LIST_FN](#)
[ZOLTAN_NUM_EDGES_MULTI_FN](#) or [ZOLTAN_NUM_EDGES_FN](#)
[ZOLTAN_EDGE_LIST_MULTI_FN](#) or [ZOLTAN_EDGE_LIST_FN](#)

Data Services and Utilities

Within Zoltan, several utilities are provided to simplify both application development and development of new algorithms in the library. They are separate from the Zoltan library so that applications can use them independently of Zoltan, if desired. They are compiled separately from Zoltan and can be archived in separate libraries. Instructions for [building the utilities](#) and [applications](#) using them are included below; individual library names are listed in the following documentation for each package.

The packages available are listed below.

- [Memory Management Utilities](#)
- [Unstructured Communication Utilities](#)
- [Distributed Directory Utility](#)

Building Utilities

The utilities provided with Zoltan have their own makefiles and can be built separately from Zoltan. If the user [builds the Zoltan library](#), the utility libraries are built automatically and copied to the appropriate *Zoltan/Obj_<platform>* directory, where *<platform>* is specified through the [ZOLTAN_ARCH environment variable](#). Zoltan and the utilities share the [Utilities/Config/Config.<platform>](#) files specifying compilation paths for various architectures. If, however, a user wishes to use these utilities without using Zoltan, he must build the libraries separately.

The structure and use of makefiles for the utilities are similar to [Zoltan's makefiles](#); a top-level makefile includes rules for building each utility's library. Object files and the utility libraries are stored in a subdirectory *Obj_<platform>*, where *<platform>* is a target architecture supported with a [Utilities/Config/Config.<platform>](#) file. The command for compiling a particular utility follows:

```
gmake ZOLTAN_ARCH=<platform> <library_name>
```

where *<library_name>* is the name of the utility library, and *<platform>* is the target architecture (corresponding to [Utilities/Config/Config.<platform>](#)). The *<library_name>* for each utility is included in the following documentation for the utilities.

Building Applications

The utilities are designed so that they can easily be used separately from Zoltan in applications. To enable type-checking of arguments, the function-prototypes file for a utility should be included in all application source code files that directly access the utility. The application must also link with the appropriate utility library (and any other libraries on which the utility depends). Library and function-prototype file names for each utility are listed in the following documentation for the utilities.

Memory Management Utilities

This package consists of wrappers around the standard C memory allocation and deallocation routines which add error-checking and [debugging capabilities](#). These routines are packaged separately from Zoltan to allow their independent use in other applications. A Fortran90 interface is not yet available. C++ programmers can include the header file "zoltan_mem.h" and use the C functions. This header file, and in fact all of Zoltan's C language header files, are surrounded by an **extern "C" {}** declaration to prevent name mangling when compiled with a C++ compiler.

Source code location:	<i>Utilities/Memory</i>
Function prototypes file:	<i>Utilities/Memory/zoltan_mem.h or include/zoltan_mem.h</i>
Library name:	libzoltan_mem.a
Other libraries used by this library:	libmpi.a. (See note below.)
Routines:	

- [Zoltan_Array_Alloc](#): Allocates arrays of dimension $n, n=0,1,...,4$
- [Zoltan_Malloc](#): Wrapper for system malloc.
- [Zoltan_Calloc](#): Wrapper for system calloc.
- [Zoltan_Realloc](#): Wrapper for system realloc.
- [Zoltan_Free](#): Frees memory and sets the pointer to NULL.
- [Zoltan_Memory_Debug](#): Sets the debug level used by the memory utilities; see the [description](#) below.
- [Zoltan_Memory_Stats](#): Prints [memory debugging](#) statistics, such as memory leak information.
- [Zoltan_Memory_Usage](#): Returns user-specified information about memory usage (i.e. maximum memory used, total memory currently allocated).
- [Zoltan_Memory_Reset](#): Sets the memory usage total specified by the user (i.e. maximum memory used, total memory currently allocated) back to zero.

Use in Zoltan:

The memory management utility routines are used extensively in Zoltan and in some individual algorithms. Zoltan developers use these routines directly for most memory management, taking advantage of the error checking and [debugging capabilities](#) of the library.

Rather than call [Zoltan_Memory_Debug](#) directly, applications using Zoltan can set the [DEBUG_MEMORY](#) parameter used by this utility through calls to [Zoltan_Set_Param](#).

Note on MPI usage:

MPI is used only to obtain the processor number (through a call to MPI_Comm_rank) for print statements and error messages. If an application does not link with MPI, the memory utilities should be compiled with -DZOLTAN_NO_MPI; all output will then appear to be from processor zero, even if it is actually from other processors.

```
double *Zoltan_Array_Alloc(char * file, int line, int n, int d1, int d2, ..., int dn, int size);
```

The **Zoltan_Array_Alloc** routine dynamically allocates an array of dimension $n, n = 0, 1, ..., 4$ with size ($d1 \times d2 \times ... \times dn$). It is intended to be used for 2, 3 and 4 dimensional arrays; [Zoltan_Malloc](#) should be used for the simpler cases. The memory allocated by **Zoltan_Array_Alloc** is contiguous, and can be freed by a single call to [Zoltan_Free](#).

Arguments:

<i>file</i>	A string containing the name of the file calling the function. The <code>__FILE__</code> macro can be passed as this argument. This argument is useful for debugging memory allocation problems.
<i>line</i>	The line number within <i>file</i> of the call to the function. The <code>__LINE__</code> macro can be passed as this argument. This argument is useful for debugging memory allocation problems.
<i>n</i>	The number of dimensions in the array to be allocated. Valid values are 0, 1, 2, 3, or 4.
<i>d1, d2, ..., dn</i>	The size of each dimension to be allocated. One argument is included for each dimension.
<i>size</i>	The size (in bytes) of the data objects to be stored in the array.

Returned Value:

double *	A pointer to the starting address of the <i>n</i> -dimensional array, or NULL if the allocation fails.
----------	--------------------------------------------------------------------------------------------------------

Example:

```
int ** x = (int **) Zoltan_Array_Alloc ( __FILE__ , __LINE__ , 2, 5, 6, sizeof (int));  
Allocates a two-dimensional, 5x6-element array of integers.
```

```
double *Zoltan_Malloc(size_t n, char *file , int line);
```

The **Zoltan_Malloc** function is a wrapper around the standard C malloc routine. It allocates a block of memory of size *n* bytes. The principle advantage of using the wrapper is that it allows memory leaks to be tracked via the `DEBUG_MEMORY` variable (set in [Zoltan_Memory_Debug](#)).

A macro **ZOLTAN_MALLOC** is defined in *zoltan_mem.h*. It takes the argument *n*, and adds the `__FILE__` and `__LINE__` macros to the argument list of the **Zoltan_Malloc** call:

```
#define ZOLTAN_MALLOC(n) Zoltan_Malloc((n), __FILE__, __LINE__)
```

Using this macro, the developer gains the file and line debugging information without having to type file and line information in each memory allocation call.

Arguments:

<i>n</i>	The size (in bytes) of the memory-allocation request.
<i>file</i>	A string containing the name of the file calling the function. The <code>__FILE__</code> macro can be passed as this argument. This argument is useful for debugging memory allocation problems.
<i>line</i>	The line number within <i>file</i> of the call to the function. The <code>__LINE__</code> macro can be passed as this argument. This argument is useful for debugging memory allocation problems.

Returned Value:

double *	A pointer to the starting address of memory allocated. NULL is returned if <i>n</i> = 0 or the routine is unsuccessful.
----------	-------------------------------------------------------------------------------------------------------------------------

Example:

```
struct Zoltan_Struct *b = (struct Zoltan_Struct *) ZOLTAN_MALLOC(sizeof(struct  
Zoltan_Struct));  
Allocates memory for one Zoltan_Struct data structure.
```

```
double *Zoltan_Calloc(size_t num, size_t size, char *file, int line);
```

The **Zoltan_Calloc** function is a wrapper around the standard C calloc routine. It allocates a block of memory of size

num * *size* bytes and initializes the memory to zeros. The principle advantage of using the wrapper is that it allows memory leaks to be tracked via the `DEBUG_MEMORY` variable (set in [Zoltan Set Memory Debug](#)).

A macro **ZOLTAN_CALLOC** is defined in *zoltan_mem.h*. It takes the arguments *num* and *size*, and adds the `__FILE__` and `__LINE__` macros to the argument list of the **Zoltan_Calloc** call:

```
#define ZOLTAN_CALLOC(num, size) Zoltan_Calloc((num), (size), __FILE__, __LINE__)
```

Using this macro, the developer gains the file and line debugging information without having to type file and line information in each memory allocation call.

Arguments:

<i>num</i>	The number of elements of the following <i>size</i> to allocate.
<i>size</i>	The size of each element. Hence, the total allocation is <i>num</i> * <i>size</i> bytes.
<i>file</i>	A string containing the name of the file calling the function. The <code>__FILE__</code> macro can be passed as this argument. This argument is useful for debugging memory allocation problems.
<i>line</i>	The line number within <i>file</i> of the call to the function. The <code>__LINE__</code> macro can be passed as this argument. This argument is useful for debugging memory allocation problems.

Returned Value:

double *	A pointer to the starting address of memory allocated. NULL is returned if <i>n</i> = 0 or the routine is unsuccessful.
----------	-------------------------------------------------------------------------------------------------------------------------

Example:

```
int *b = (int *) ZOLTAN_CALLOC( 10, sizeof(int));  
Allocates memory for 10 integers and initializes the memory to zeros.
```

```
double *Zoltan_Realloc(void *ptr, size_t n, char *file, int line);
```

The **Zoltan_Realloc** function is a "safe" version of `realloc`. It changes the size of the object pointed to by *ptr* to *n* bytes. The contents of *ptr* are unchanged up to a minimum of the old and new sizes. Error tests ensuring that *n* is a positive number and that space is available to be allocated are performed.

A macro **ZOLTAN_REALLOC** is defined in *zoltan_mem.h*. It takes the arguments *ptr* and *n*, and adds the `__FILE__` and `__LINE__` macros to the argument list of the **Zoltan_Realloc** call:

```
#define ZOLTAN_REALLOC(ptr, n) Zoltan_Realloc((ptr), (n), __FILE__, __LINE__)
```

Using this macro, the developer gains the file and line debugging information without having to type file and line information in each memory allocation call.

Arguments:

<i>ptr</i>	Pointer to allocated memory to be re-sized.
<i>n</i>	The size (in bytes) of the memory-allocation request.
<i>file</i>	A string containing the name of the file calling the function. The <code>__FILE__</code> macro can be passed as this argument. This argument is useful for debugging memory allocation problems.
<i>line</i>	The line number within <i>file</i> of the call to the function. The <code>__LINE__</code> macro can be passed as this argument. This argument is useful for debugging memory allocation problems.

Returned Value:

double *	A pointer to the starting address of memory allocated. If the routine is unsuccessful, NULL is returned and <i>*ptr</i> is unchanged.
----------	---------------------------------------------------------------------------------------------------------------------------------------

Example:

```
int n = sizeof(struct Zoltan_Struct);
int *b = (int *) ZOLTAN\_MALLOC (n);
b = (int *) ZOLTAN\_REALLOC (b, 2*n);
Reallocates memory for b from length n to length 2*n.
```

```
void Zoltan_Free(void **ptr, char * file , int line);
```

The **Zoltan_Free** function calls the system's "free" function for the memory pointed to by **ptr*. Note that the argument to this routine has an extra level of indirection when compared to the standard C "free" call. This allows the pointer being freed to be set to NULL, which can help find errors in which a pointer is used after it is deallocated. Error checking is performed to prevent attempts to free NULL pointers. When **Zoltan_Free** is used with the `DEBUG_MEMORY` options (set in [Zoltan_Memory_Debug](#)), it can help identify memory leaks.

A macro **ZOLTAN_FREE** is defined in *zoltan_mem.h*. It takes the argument *ptr*, and adds the `__FILE__` and `__LINE__` macros to the argument list of the **Zoltan_Free** call:

```
#define ZOLTAN_FREE(ptr) Zoltan_Free((void **)(ptr), __FILE__, __LINE__)
```

Using this macro, the developer gains the file and line debugging information without having to type file and line information in each memory allocation call.

Arguments:

ptr Address of a pointer to the memory to be freed. Upon return, *ptr* is set to NULL.

Example:

```
ZOLTAN_FREE(& x);
Frees memory associated with the variable x; upon return, x is NULL.
```

Debugging Memory Errors

One important reason to use the memory-management utilities' wrappers around the system memory routines is to facilitate debugging of memory problems. Various amounts of information can about memory allocation and deallocation are stored, depending on the debug level set through a call to [Zoltan_Memory_Debug](#). This information is printed either when an error or warning occurs, or when [Zoltan_Memory_Stats](#) is called. We have found values of one and two to be very helpful in our development efforts. The routine [Zoltan_Memory_Usage](#) can be called to return user-specified information about memory utilization to the user's program, and [Zoltan_Memory_Reset](#) can be called to set totals back to zero.

```
void Zoltan_Memory_Debug(int new_level);
```

The **Zoltan_Memory_Debug** function sets the level of memory debugging to be used.

Arguments:

new_level Integer indicating the amount of debugging to use. Valid options include:

- 0 -- No debugging.
- 1 -- The number of calls to [Zoltan_Malloc](#) and [Zoltan_Free](#) are tallied, and can

be printed by a call to [Zoltan_Memory_Stats](#).
2 -- A list of all calls to [Zoltan_Malloc](#) which have not yet been freed is kept. This list is printed by [Zoltan_Memory_Stats](#) (useful for detecting memory leaks). Any calls to [Zoltan_Free](#) with addresses not in this list trigger warning messages. (Note that allocations that occurred prior to setting the debug level to 2 will not be in this list and thus can generate spurious warnings.)
3 -- Information about each allocation is printed as it happens.

Default:

Memory debug level is 1.

void **Zoltan_Memory_Stats**();

The **Zoltan_Memory_Stats** function prints information about memory allocation and deallocation. The amount of information printed is determined by the debug level set through a call to [Zoltan_Memory_Debug](#).

Arguments:

None.

size_t **Zoltan_Memory_Usage**(int *type*);

The **Zoltan_Memory_Usage** function returns information about memory utilization. The memory debug level (set through a call to [Zoltan_Set_Memory_Debug](#)) must be at least 2 for this function to return non-zero values.

Arguments:

type Integer to request type of information required. These integers are defined in *zoltan_mem.h*. Valid options include:

- ZOLTAN_MEM_STAT_TOTAL -- The function will return the current total memory allocated via Zoltan's memory allocation routines.
- ZOLTAN_MEM_STAT_MAXIMUM -- The function will return the maximum total memory allocated via Zoltan's memory allocation routines up to this point.

Default:

type = ZOLTAN_MEM_STAT_MAXIMUM

Returned Value:

int The number in bytes of the specific requested memory statistic.

Example:

total = **Zoltan_Memory_Usage** (ZOLTAN_MEM_STAT_TOTAL);

void **Zoltan_Memory_Reset**(int *type*);

The **Zoltan_Memory_Reset** function sets the specified count to zero.

Arguments:

type Integer to specify the type of information to be reset . These integers are defined in *zoltan_mem.h*. Valid options include:

ZOLTAN_MEM_STAT_TOTAL -- The function will set the count of total memory allocated via Zoltan's memory allocation routines to zero.
ZOLTAN_MEM_STAT_MAXIMUM -- The function will set the count of maximum total memory allocated via Zoltan's memory allocation routines back to zero.

Default:

type = ZOLTAN_MEM_STAT_MAXIMUM

Example:

Zoltan_Memory_Reset (*ZOLTAN_MEM_STAT_TOTAL*);

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Unstructured Communication Utilities

The unstructured communication package provides a simple interface for doing complicated patterns of point-to-point communication, such as those associated with data remapping. This package consists of a few simple functions which create or modify communication *plans*, perform communication, and destroy communication plans upon completion. The package is descended from software first developed by Steve Plimpton and Bruce Hendrickson, and has proved useful in a variety of different applications. For this reason, it is maintained as a separate library and can be used independently from Zoltan.

In a prototypical usage of the unstructured communication package each processor has some objects to send to other processors, but no processor knows what messages it will receive. A call to [Zoltan_Comm_Create](#) produces a data structure called a communication *plan* which encapsulates the basic information about the communication operation. The plan does not know anything about the types of objects being transferred, only the number of them. So the same plan can be used repeatedly to transfer different types of data as long as the number of objects in the transfers remains the same. The actual size of objects isn't specified until the call to [Zoltan_Comm_Do](#) which performs the data transfer.

The plan which is produced by [Zoltan_Comm_Create](#) assumes that all the objects are of the same size. If this is not true, then a call to [Zoltan_Comm_Resize](#) can specify the actual size of each object, and the plan is augmented appropriately. [Zoltan_Comm_Resize](#) can be invoked repeatedly on the same plan to specify varying sizes for different data transfer operations.

Although a friendlier interface may be added in the future, for now the data to be sent must be passed to [Zoltan_Comm_Do](#) as a packed buffer in which the objects are stored consecutively. This probably requires the application to pull the data out of native data structures and place in into the buffer. The destination of each object is specified by the *proclist* argument to [Zoltan_Comm_Create](#). Some flexibility is supported by allowing *proclist* to contain negative values, indicating that the corresponding objects are not to be sent. The communication operations allow for any object to be sent to any destination processor. However, if the objects are grouped in such a way that all those being sent to a particular processor are consecutive, the time and memory of an additional copy is avoided.

Function [Zoltan_Comm_Do_Reverse](#) reverses the communication plan to send back messages to the originators.

To allow overlap between communication and processing, POST and WAIT variants of [Zoltan_Comm_Do](#) and [Zoltan_Comm_Do_Reverse](#) are provided. Communication is initiated by the POST function ([Zoltan_Comm_Do_Post](#) or [Zoltan_Comm_Do_Reverse_Post](#)); incoming messages are posted and outgoing messages are sent. Then the user can continue processing. After the processing is complete, the corresponding WAIT function ([Zoltan_Comm_Do_Wait](#) or [Zoltan_Comm_Do_Reverse_Wait](#)) is called to wait for all incoming messages to be received. For convenience, these functions use the same calling arguments as [Zoltan_Comm_Do](#) and [Zoltan_Comm_Do_Reverse](#).

All the functions in the unstructured communication library return integer [error codes](#) identical to those used by Zoltan.

The C++ interface to the unstructured communication utility is found in the `zoltan_comm_cpp.h` header file which defines the `Zoltan_Comm` class.

A Fortran90 interface is not yet available.

Source code location:	Utilities/Communication
C Function prototypes file:	Utilities/Communication/zoltan_comm.h
C++ class definition:	Utilities/Communication/zoltan_comm_cpp.h

Library name: libzoltan_comm.a
Other libraries used by this library: libmpi.a, [libzoltan_mem.a](#).
High Level Routines:

[Zoltan_Comm_Create](#): computes a communication plan for sending objects to destination processors.

[Zoltan_Comm_Do](#): uses a communication plan to send data objects to destination processors. The POST and WAIT variants are

[Zoltan_Comm_Do_Post](#) and

[Zoltan_Comm_Do_Wait](#).

[Zoltan_Comm_Do_Reverse](#): performs the reverse (opposite) communication of **[Zoltan_Comm_Do](#)**.

The POST and WAIT variants are

[Zoltan_Comm_Do_Reverse_Post](#) and

[Zoltan_Comm_Do_Reverse_Wait](#).

[Zoltan_Comm_Resize](#): augments the plan to allow objects to be of variable sizes.

[Zoltan_Comm_Copy](#): create a new communication plan and copy an existing one to it.

[Zoltan_Comm_Copy_To](#): copy one existing communication plan to another.

[Zoltan_Comm_Destroy](#): free memory associated with a communication plan.

Low Level Routines:

[Zoltan_Comm_Exchange_Sizes](#): updates the sizes of the messages each processor will receive.

[Zoltan_Comm_Invert_Map](#): given a set of messages each processor wants to send, determines the set of messages each processor needs to receive.

[Zoltan_Comm_Sort_Ints](#): sorts an array of integer values.

[Zoltan_Comm_Info](#): returns information about a communication plan.

[Zoltan_Comm_Invert_Plan](#): given a communication plan, converts the plan into a plan for the reverse communication.

Use in Zoltan:

The Zoltan library uses the unstructured communication package in its migration tools and in some of the load-balancing algorithms. For example, in **[Zoltan_Migrate](#)**, **[Zoltan_Comm_Create](#)** is used to develop a communication map for sending objects to be exported to their new destination processors. The sizes of the exported objects are obtained and the communication map is augmented with a call to **[Zoltan_Comm_Resize](#)**. The data for the objects is packed into a communication buffer and sent to the other processors through a call to **[Zoltan_Comm_Do](#)**. After the received objects are unpacked, the communication plan is no longer needed, and it is deallocated by a call to **[Zoltan_Comm_Destroy](#)**. Zoltan developers use the package whenever possible, as improvements made to the package (for example, support for heterogeneous architectures) automatically propagate to the algorithms.

C:

```
int Zoltan_Comm_Create( struct Zoltan_Comm_Obj **plan, int nsend, int *proclist, MPI_Comm comm, int tag, int *nreturn);
```

C++:

```
Zoltan_Comm( const int & nsend, int *proclist, const MPI_Comm & comm, const int & tag, int *nreturn);
```

or

```
Zoltan_Comm();
```

```
Zoltan_Comm::Create( const int & nsend, int *proclist, const MPI_Comm & comm, const int & tag, int *nreturn);
```

The **[Zoltan_Comm_Create](#)** function sets up the communication plan in the unstructured communication package. Its input is a count of objects to be sent to other processors, a list of the processors to which the objects should be sent (repetitions are allowed), and an MPI communicator and tag. It allocates and builds a communication plan that

describes to which processors data will be sent and from which processors data will be received. It also computes the amount of data to be sent to and received from each processor. It returns the number of objects to be received by the processor and a pointer to the communication plan it created. The communication plan is then used by calls to [Zoltan_Comm_Do](#) to perform the actual communication.

Arguments:

<i>plan</i>	A pointer to the communication plan created by Zoltan_Comm_Create .
<i>nsend</i>	The number of objects to be sent to other processors.
<i>proclist</i>	An array of size <i>nsend</i> of destination processor numbers for each of the objects to be sent.
<i>comm</i>	The MPI communicator for the unstructured communication.
<i>tag</i>	A tag for MPI communication.
<i>nreturn</i>	Upon return, the number of objects to be received by the processor.

Returned Value:

int	Error code.
-----	-------------

In the C++ interface to the communication utility, the communication plan is represented by a **Zoltan_Comm** object. It is created when the **Zoltan_Comm** constructor executes. There are two constructors. The first one listed above uses parameters to initialize the plan. The second constructor does not, but the plan can subsequently be initialized with a call to **Zoltan_Comm::Create()**.

C:

```
int Zoltan_Comm_Do(struct Zoltan_Comm_Obj *plan, int tag, char *send_data, int nbytes, char *recvbuf);
int Zoltan_Comm_Do_Post(struct Zoltan_Comm_Obj *plan, int tag, char *send_data, int nbytes, char *recvbuf);
int Zoltan_Comm_Do_Wait(struct Zoltan_Comm_Obj *plan, int tag, char *send_data, int nbytes, char *recvbuf);
```

C++:

```
int Zoltan_Comm::Do(const int & tag, char *send_data, const int & nbytes, char *recvbuf);
int Zoltan_Comm::Do_Post(const int & tag, char *send_data, const int & nbytes, char *recvbuf);
int Zoltan_Comm::Do_Wait(const int & tag, char *send_data, const int & nbytes, char *recvbuf);
```

The **Zoltan_Comm_Do** function performs the communication described in a communication plan built by [Zoltan_Comm_Create](#). Using the plan, it takes a buffer of object data to be sent and the size (in bytes) of each object's data in that buffer and sends the data to other processors. **Zoltan_Comm_Do** also receives object data from other processors and stores it in a receive buffer. The receive buffer must be allocated by the code calling **Zoltan_Comm_Do** using the number of received objects returned by [Zoltan_Comm_Create](#) or [Zoltan_Comm_Resize](#). If the objects have variable sizes, then [Zoltan_Comm_Resize](#) must be called before **Zoltan_Comm_Do**.

Arguments:

<i>plan</i>	A pointer to a communication plan built by Zoltan_Comm_Create .
<i>tag</i>	An MPI message tag.
<i>send_data</i>	A buffer filled with object data to be sent to other processors.
<i>nbytes</i>	The size (in bytes) of the data for one object, or the scale factor if the objects have variable sizes. (See Zoltan_Comm_Resize for more details.)
<i>recvbuf</i>	Upon return, a buffer filled with object data received from other processors.

Returned Value:

int	Error code.
-----	-------------

```
C:
int Zoltan_Comm_Do_Reverse( struct Zoltan_Comm_Obj *plan, int tag, char *send_data, int nbytes, int *sizes, char
*recvbuf);
int Zoltan_Comm_Do_Reverse_Post( struct Zoltan_Comm_Obj *plan, int tag, char *send_data, int nbytes, int *sizes,
char *recvbuf);
int Zoltan_Comm_Do_Reverse_Wait( struct Zoltan_Comm_Obj *plan, int tag, char *send_data, int nbytes, int
*sizes, char *recvbuf);
C++:
int Zoltan_Comm::Do_Reverse( const int & tag, char *send_data, const int & nbytes, int *sizes, char *recvbuf);
int Zoltan_Comm::Do_Reverse_Post( const int & tag, char *send_data, const int & nbytes, int *sizes, char
*recvbuf);
int Zoltan_Comm::Do_Reverse_Wait( const int & tag, char *send_data, const int & nbytes, int *sizes, char
*recvbuf);
```

The **Zoltan_Comm_Do_Reverse** function performs communication based on a communication plan built by [Zoltan_Comm_Create](#). But unlike [Zoltan_Comm_Do](#), this routine performs the reverse of the communication pattern. Specifically, all sends in the plan are treated as receives and vice versa. **Zoltan_Comm_Do_Reverse** is particularly well suited to return updated data objects to their originating processors when the objects were initially transferred via [Zoltan_Comm_Do](#).

Arguments:

<i>plan</i>	A pointer to a communication plan built by Zoltan_Comm_Create .
<i>tag</i>	An MPI message tag to be used by this routine.
<i>send_data</i>	A buffer filled with object data to be sent to other processors.
<i>nbytes</i>	The size (in bytes) of the data associated with an object, or the scale factor if the objects have variable sizes.
<i>sizes</i>	If not NULL, this input array specifies the size of all the data objects being transferred. This argument is passed directly to Zoltan_Comm_Resize . This array has length equal to the <i>nsend</i> value passed to Zoltan_Comm_Create . But note that for Zoltan_Comm_Do_Reverse this array describes the sizes of the values being received, not sent.
<i>recvbuf</i>	Upon return, a buffer filled with object data received from other processors.

Returned Value:

int	Error code.
-----	-------------

```
C:
int Zoltan_Comm_Resize( struct Zoltan_Comm_Obj *plan, int *sizes, int tag , int *total_recv_size);
C++:
int Zoltan_Comm::Resize( int *sizes, const int & tag , int *total_recv_size);
```

If the objects being communicated are of variable sizes, then the plan produced by [Zoltan_Comm_Create](#) is incomplete. This routine allows the plan to be augmented to allow for variable sizes. **Zoltan_Comm_Resize** can be invoked repeatedly on the same plan to specify different object sizes associated with different data transfers.

Arguments:

<i>plan</i>	A communication plan built by Zoltan_Comm_Create .
<i>sizes</i>	An input array of length equal to the <i>nsend</i> argument in the call to Zoltan_Comm_Create which generated the <i>plan</i> . Each entry in the array is the size of the corresponding object to be sent. If <i>sizes</i> is <i>NULL</i> (on all processors), the objects are considered to be the same size. Note that the true size of a message will be scaled by the <i>nbytes</i> argument to Zoltan_Comm_Do .

<i>tag</i>	A message tag to be used for communication within this routine, based upon the communicator in <i>plan</i> .
<i>total_recv_size</i>	Sum of the sizes of the incoming messages. To get the actual size (in bytes), you need to scale by the <i>nbytes</i> argument to Zoltan_Comm_Do .
Returned Value:	
int	Error code.

```
C: struct Zoltan_Comm_Obj *Zoltan_Comm_Copy( struct Zoltan_Comm_Obj *plan);
C++: Zoltan_Comm(const Zoltan_Comm &plan);
```

Zoltan_Comm_Copy creates a new *Zoltan_Comm_Obj* structure and copies the existing *plan* to it. The corresponding C++ method is the **Zoltan_Comm** copy constructor.

Arguments:	
<i>plan</i>	A pointer to the communication plan to be copied to the new <i>Zoltan_Comm_Obj</i> structure.
Returned Value:	
struct Zoltan_Comm_Obj *	the newly created plan, or NULL on error.

```
C: int Zoltan_Comm_Copy_To( struct Zoltan_Comm_Obj **to, struct Zoltan_Comm_Obj *from);
C++: Zoltan_Comm & operator= (const Zoltan_Comm &plan);
```

Zoltan_Comm_Copy_To copies one existing communication plan to another. The corresponding C++ method is the **Zoltan_Comm** copy operator.

Arguments:	
<i>to</i>	A pointer to a pointer to the communication plan that will be copied to. We destroy the plan first, and set the pointer to the plan to NULL, before proceeding with the copy.
<i>from</i>	A pointer the communication plan that we will make a copy of.
Returned Value:	
int	Error code

```
C: int Zoltan_Comm_Destroy( struct Zoltan_Comm_Obj **plan);
C++: ~Zoltan_Comm();
```

The **Zoltan_Comm_Destroy** function frees all memory associated with a communication plan created by [Zoltan_Comm_Create](#). The C++ **Zoltan_Comm** object does not have an explicit **Destroy** method. It is deallocated when its destructor is called.

Arguments:	
<i>plan</i>	A pointer to a communication plan built by Zoltan_Comm_Create . Upon return, <i>plan</i> is set to NULL.
Returned Value:	
int	Error code.

```
C:
int Zoltan_Comm_Exchange_Sizes( int *sizes_to, int *procs_to, int nsends, int self_msg, int *sizes_from, int
*procs_from, int nrecvs, int *total_recv_size, int my_proc, int tag, MPI_Comm comm );

C++:
static int Zoltan_Comm::Exchange_Sizes( int *sizes_to, int *procs_to, const int & nsends, const int & self_msg, int
*sizes_from, int *procs_from, const int & nrecvs, int *total_recv_size, const int & my_proc, const int & tag, const
MPI_Comm & comm );
```

This routine is used by [Zoltan_Comm_Resize](#) to update the sizes of the messages each processor is expecting to receive. The processors already know who will send them messages, but if variable sized objects are being communicated, then the sizes of the messages are recomputed and exchanged via this routine.

Arguments:

<i>sizes_to</i>	Input array with the size of each message to be sent. Note that the actual number of bytes in the message is the product of this value and the <i>nbytes</i> argument to Zoltan_Comm_Do .
<i>procs_to</i>	Input array with the destination processor for each of the messages to be sent.
<i>nsends</i>	Input argument with the number of messages to be sent. (Length of the <i>procs_to</i> array.)
<i>self_msg</i>	Input argument indicating whether a processor has data for itself (=1) or not (=0) within the <i>procs_to</i> and <i>lengths_to</i> arrays.
<i>sizes_from</i>	Returned array with the size of each message that will be received. Note that the actual number of bytes in the message is the product of this value and the <i>nbytes</i> argument to Zoltan_Comm_Do .
<i>procs_from</i>	Returned array of processors from which data will be received.
<i>nrecvs</i>	Returned value with number of messages to be received. (length of <i>procs_from</i> array.)
<i>total_recv_size</i>	The total size of all the messages to be received. As above, the actual number of bytes will be scaled by the <i>nbytes</i> argument to Zoltan_Comm_Do .
<i>my_proc</i>	The processor's ID in the <i>comm</i> communicator.
<i>tag</i>	A message tag which can be used by this routine.
<i>comm</i>	MPI Communicator for the processor numbering in the <i>procs</i> arrays.

Returned Value:

int	Error code.
-----	-------------

```
C:
int Zoltan_Comm_Invert_Map( int *lengths_to, int *procs_to, int nsends, int self_msg, int ** lengths_from, int **
procs_from, int * nrecvs, int my_proc, int nprocs, int out_of_mem, int tag, MPI_Comm comm );

C++:
static int Zoltan_Comm::Invert_Map( int *lengths_to, int *procs_to, const int & nsends, const int & self_msg, int *
& lengths_from, int * & procs_from, int & nrecvs, const int & my_proc, const int & nprocs, const int & out_of_mem,
const int & tag, const MPI_Comm & comm );
```

The **Zoltan_Comm_Invert_Map** function is a low level communication routine. It is useful when a processor knows to whom it needs to send data, but not from whom it needs to receive data. Each processor provides to this routine a set of lengths and destinations for the messages it wants to send. The routine then returns the set of lengths and origins for the messages a processor will receive. Note that by inverting the interpretation of *to* and *from* in these arguments, the routine can be used to do the opposite: knowing how much data to receive and from which processors, it can compute how much data to send and to which processors.

Arguments:

<i>lengths_to</i>	Input array with the number of values in each of the messages to be sent. Note that the actual size of each value is not specified until the Zoltan_Comm_Do routine is invoked.
<i>procs_to</i>	Input array with the destination processor for each of the messages to be sent.
<i>nsends</i>	Input argument with the number of messages to be sent. (Length of the <i>lengths_to</i> and <i>procs_to</i> arrays.)
<i>self_msg</i>	Input argument indicating whether a processor has data for itself (=1) or not (=0) within the <i>procs_to</i> and <i>lengths_to</i> arrays.
<i>lengths_from</i>	Returned array with lengths of messages to be received.
<i>procs_from</i>	Returned array of processors from which data will be received.
<i>nrecvs</i>	Returned value with number of messages to be received (lengths of <i>lengths_from</i> and <i>procs_from</i> arrays).
<i>my_proc</i>	The processor's ID in the <i>comm</i> communicator.
<i>nprocs</i>	Number of processors in the <i>comm</i> communicator.
<i>out_of_mem</i>	Since it has a barrier operation, this routine is a convenient time to tell all the processors that one of them is out of memory. This input argument is 0 if the processor is OK, and 1 if the processor has failed in a malloc call. All the processors will return with a code of COMM_MEMERR if any of them is out of memory.
<i>tag</i>	A message tag which can be used by this routine.
<i>comm</i>	MPI Communicator for the processor numbering in the <i>procs</i> arrays.
Returned Value:	
int	Error code.

int **Zoltan_Comm_Sort_Ints**(int *vals_sort, int *vals_other, int nvals);

As its name suggests, the **Zoltan_Comm_Sort_Ints** function sorts a set of integers via the quicksort algorithm. The integers are reordered from lowest to highest, and a second array of integers is reordered in the same fashion. This second array can be used to return the permutation associated with the sort operation. There is no C++ interface to this function. You can use the C function instead.

Arguments:

<i>vals_sort</i>	The array of integers to be sorted. This array is permuted into sorted order.
<i>vals_other</i>	Another array of integers which is permuted identically to <i>vals_sort</i> .
<i>nvals</i>	The number of values in the two integer arrays.

Returned Value:

int	Error code.
-----	-------------

C:

int **Zoltan_Comm_Info**(struct Zoltan_Comm_Obj *plan, int *nsends, int *send_procs, int *send_lengths, int *send_nvals, int *send_max_size, int *send_list, int *nrecvs, int *recv_procs, int *recv_lengths, int *recv_nvals, int *recv_total_size, int *recv_list, int *self_msg);

C++:

int **Zoltan_Comm::Info**(int *nsends, int *send_procs, int *send_lengths, int *send_nvals, int *send_max_size, int *send_list, int *nrecvs, int *recv_procs, int *recv_lengths, int *recv_nvals, int *recv_total_size, int *recv_list, int *self_msg) const;

Zoltan_Comm_Info returns information about a communication plan. All arguments, except the *plan* itself, may be NULL; values are returned only for non-NULL arguments.

Arguments:

<i>plan</i>	Communication data structure created by Zoltan_Comm_Create .
<i>nsends</i>	Upon return, the number of processors to which messages are sent; does not include self-messages.
<i>send_procs</i>	Upon return, a list of processors to which messages are sent; self-messages are included.
<i>send_lengths</i>	Upon return, the number of values to be sent to each processor in <i>send_procs</i> .
<i>send_nvals</i>	Upon return, the total number of values to send.
<i>send_max_size</i>	Upon return, the maximum size of a message to be sent; does not include self-messages.
<i>send_list</i>	Upon return, the processor assignment of each value to be sent.
<i>nrecvs</i>	Upon return, the number of processors from which to receive messages; does not include self-messages.
<i>recv_procs</i>	Upon return, a list of processors from which messages are received; includes self-messages.
<i>recv_lengths</i>	Upon return, the number of values to be received from each processor in <i>recv_procs</i> .
<i>recv_nvals</i>	Upon return, the total number of values to receive.
<i>recv_total_size</i>	Upon return, the total size of items to be received.
<i>recv_list</i>	Upon return, the processor assignments of each value to be received.
<i>self_msg</i>	Upon return, the number of self-messages.

Returned Value:

int	Error code.
-----	-------------

```
C: int Zoltan_Comm_Invert_Plan( struct Zoltan_Comm_Obj **plan );
C++: int Zoltan_Comm::Invert_Plan();
```

Given a communication plan, **Zoltan_Comm_Invert_Plan** alters the plan to make it the plan for the reverse communication. Information in the input plan is replaced by information for the reverse-communication plan. All receives in the reverse-communication plan are blocked; thus, using the inverted plan does not produce the same results as [Zoltan_Comm_Do_Reverse](#). If an error occurs within **Zoltan_Comm_Invert_Plan**, the original plan is returned unaltered.

Arguments:

<i>plan</i>	Communication data structure created by Zoltan_Comm_Create ; the contents of this plan are irretrievably modified by Zoltan_Comm_Invert_Plan .
-------------	-----------------------------------------------------------------------------------------------------------------------------------------------------------------------

Returned Value:

int	Error code.
-----	-------------

Distributed Directory Utility

A distributed directory may be viewed as a distributed hash table pointing to the information stored in the directory. An application may use this directory utility to manage its objects' locations after data migrations or to make information globally accessible. A distributed directory balances the load (in terms of memory and processing time) and avoids the bottle neck of a centralized directory design.

This distributed directory module may be used alone or in conjunction with Zoltan's load balancing capability and memory and communication services. The user should note that external names (subroutines, etc.) prefaced by `Zoltan_DD_` are reserved when using this module. Since the distributed directory uses collective communication, it is important that all processors call the same function at the same time in their processing.

The user initially creates an empty distributed directory using [Zoltan_DD_Create](#). Then each global ID (GID), which are the directory keys, together with other optional information is added to the directory using [Zoltan_DD_Update](#). The directory maintains the GID's basic information: local ID (optional), part (optional), arbitrary user data (optional), and the current data owner (optional). [Zoltan_DD_Update](#) is also called after data migration or whenever it is useful to update the information in the directory. [Zoltan_DD_Find](#) returns the directory information for a list of GIDs. A selected list of GIDs may be removed from the directory by [Zoltan_DD_Remove](#). When the user has finished using the directory, its memory is returned to the system by [Zoltan_DD_Destroy](#).

An object is known by its GID. Hashing provides very fast lookup for the information associated with a GID in a two step process. The first hash of the GID yields the processor number owning the directory entry for that GID. The directory entry owner remains constant even if the object associated with the GID migrates or changes over time. Second, a different hash algorithm on the GID looks up the associated information in directory processor's hash table. The user may optionally register their own (first) hash function to take advantage of their knowledge of their GID naming scheme and the GID's neighboring processors. See the documentation for [Zoltan_DD_Set_Hash_Fn](#) for more information. If no user hash function is registered, Zoltan's [Zoltan_Hash](#) will be used. This module's design was strongly influenced by the paper "Communication Support for Adaptive Computation" by Pinar and Hendrickson.

Some users number their GIDs by giving the first "n" GIDs to processor 0, the next "n" GIDs to processor 1, and so forth. The function [Zoltan_DD_Set_Neighbor_Hash_Fn1](#) will provide efficient directory communication when these GIDs stay close to their origin. The function `Zoltan_DD_Set_Neighbor_Hash_Fn2` allows the specification of ranges of GIDs to each processor for more flexibility. The source code for [DD_Set_Neighbor_Hash_Fn1](#) and [DD_Set_Neighbor_Hash_Fn2](#) provide examples of how a user can create their own "hash" functions taking advantage of their own GID naming convention.

The routine [Zoltan_DD_Print](#) will print the contents of the directory. The companion routine [Zoltan_DD_Stats](#) prints out a summary of the hash table size, number of linked lists, and the length of the longest linked list. This may be useful when the user creates their own hash functions.

All modules use the following response to the `debug_level`:

`debug_level=0`, Output is silent except for FATAL or MEMERR errors.

`debug_level=1`, Turns on checking for legal, but possibly wrong conditions such as updating the same directory multiple times in one update cycle.

`debug_level=5`, Adds tracing information for the routines defined below.

`debug_level=6`, Adds tracing information for all DD routines.

`debug_level=7`, Adds tracing within each routine,

`debug_level>7`, Adds information about each object when used.

Calling `DD_Stats` or `DD_Print` is automatically verbose independent of the `debug_level`.

The C++ interface to this utility is defined in the header file *zoltan_dd_cpp.h* as the class **Zoltan_DD**. A single **Zoltan_DD** object represents a distributed directory.

A Fortran90 interface is not yet available.

Source code location:	<i>Utilities/DDirectory</i>
C Function prototypes file:	<i>Utilities/DDirectory/zoltan_dd.h</i>
C++ class definition:	<i>Utilities/DDirectory/zoltan_dd_cpp.h</i>
Library name:	libzoltan_dd.a
Other libraries used by this library:	libmpi.a, libzoltan_mem.a, libzoltan_comm.a
Routines:	

- [**Zoltan_DD_Create**](#): Allocates memory and initializes the directory.
- [**Zoltan_DD_Copy**](#): Allocates a new directory structure and copies an existing one to it.
- [**Zoltan_DD_Copy_To**](#): Copies one directory structure to another.
- [**Zoltan_DD_Destroy**](#): Terminate the directory and frees its memory.
- [**Zoltan_DD_Update**](#): Adds or updates GIDs' directory information.
- [**Zoltan_DD_Find**](#): Returns GIDs' information (owner, local ID, etc.)
- [**Zoltan_DD_Remove**](#): Eliminates selected GIDs from the directory.
- [**Zoltan_DD_Stats**](#): Provides statistics about hash table & linked lists.
- [**Zoltan_DD_Print**](#): Displays the contents (GIDs, etc) of each directory.
- [**Zoltan_DD_Set Hash Fn**](#): Registers a user's optional hash function.
- [**Zoltan_DD_Set Neighbor Hash Fn1**](#): Hash function with constant number of GIDs per processor.
- [**Zoltan_DD_Set Neighbor Hash Fn2**](#): Hash function with variable number of GID's per processor.

Data Structures:

struct Zoltan_DD_Struct: State & storage used by all DD routines. Users should not modify any internal values in this structure. Users should only pass the address of this structure to the other routines in this package.

C:	<code>int Zoltan_DD_Create (struct Zoltan_DD_Struct **dd, MPI_Comm comm, int num_gid_entries, int num_lid_entries, int user_length, int table_length, int debug_level);</code>
C++:	<code>Zoltan_DD(const MPI_Comm & comm, const int & num_gid_entries, const int & num_lid_entries, const int & user_length, const int & table_length, const int & debug_level);</code> or <code>Zoltan_DD();</code> <code>Zoltan_DD::Create(const MPI_Comm & comm, const int & num_gid_entries, const int & num_lid_entries, const int & user_length, const int & table_length, const int & debug_level);</code>

Zoltan_DD_Create allocates and initializes memory for the **Zoltan_DD_Struct** structure. It must be called before any other distributed directory routines. MPI must be initialized prior to calling this routine.

The **Zoltan_DD_Struct** must be passed to all other distributed directory routines. The MPI Comm argument designates the processors used for the distributed directory. The MPI Comm argument is duplicated and stored for later use. The length of the GID, length of the LID, and the length of the optional user data (*user_length*) must be consistent for all processors.

Arguments:

<i>dd</i>	Structure maintains directory state and hash table.
<i>comm</i>	MPI comm duplicated and stored specifying directory processors.
<i>num_gid_entries</i>	Length of GID.
<i>num_lid_entries</i>	Length of local ID or zero to ignore local IDs.
<i>user_length</i>	Length of user defined data field (optional, may be zero).
<i>table_length</i>	Length of hash table (zero forces default value of 100,000 slots). For large problems, this value should be increased to approximately the number of global GIDs / number of processors (if you have enough memory) in order to improve performance.
<i>debug_level</i>	Legal values range in [0,9]. Sets the output response to various error conditions where 9 is the most verbose.

Returned Value:

int	Error code .
-----	------------------------------

ZOLTAN_FATAL is returned for MPI problems or if *num_gid_entries*, *num_lid_entries*, or *user_length* do not match globally.
ZOLTAN_MEMERR is returned if sufficient memory can not be allocated.
ZOLTAN_OK is the normal return value.

In the C++ interface, the distributed directory is represented by a **Zoltan_DD** object. It is created when the **Zoltan_DD** constructor executes. There are two constructors. The first one listed above uses parameters to initialize the distributed directory. The second constructor does not, but it can subsequently be initialized with a call to **Zoltan_DD::Create()**.

C:	struct Zoltan_DD_Struct * Zoltan_DD_Copy (struct Zoltan_DD_Struct * <i>from</i>);
C++:	Zoltan_DD (const Zoltan_DD &dd);

This routine creates a new distributed directory structure and copies an existing one to it. The corresponding routine in the C++ library is the Zoltan_DD copy constructor.

Arguments:

<i>from</i>	The existing directory structure which will be copied to the new one.
-------------	-----------------------------------------------------------------------

Returned Value:

struct Zoltan_DD_Struct *	The newly created directory structure.
------------------------------	----------------------------------------

C:	int Zoltan_DD_Copy_To (struct Zoltan_DD_Struct ** <i>to</i> , struct Zoltan_DD_Struct * <i>from</i>);
C++:	Zoltan_DD & operator =(const Zoltan_DD &dd);

This routine copies one distributed directory structure to another. The corresponding method in the C++ library is the Zoltan_DD class copy operator.

Arguments:

<i>to</i>	A pointer to a pointer to the target structure. The structure will be destroyed and the pointer set to NULL before proceeding with the copy.
<i>from</i>	A pointer to the source structure. The contents of this structure will be copied to the target structure.

Returned Value:

int [Error code](#).

C: void **Zoltan_DD_Destroy** (struct Zoltan_DD_Struct ***dd*);
C++: ~**Zoltan_DD**();

This routine frees all memory allocated for the distributed directory. No calls to any distributed directory functions using this Zoltan_DD_Struct are permitted after calling this routine. MPI is necessary for this routine only to free the previously saved MPI comm.

Arguments:

dd Directory structure to be deallocated.

Returned Value:

void NONE

There is no explicit **Destroy** method in the C++ **Zoltan_DD** class. The object is deallocated when its destructor is called.

C: int **Zoltan_DD_Update** (struct Zoltan_DD_Struct **dd*, [ZOLTAN_ID_PTR](#) *gid*, [ZOLTAN_ID_PTR](#) *lid*, [ZOLTAN_ID_PTR](#) *user*, int **part*, int *count*);
C++: int **Zoltan_DD::Update**([ZOLTAN_ID_PTR](#) *gid*, [ZOLTAN_ID_PTR](#) *lid*, [ZOLTAN_ID_PTR](#) *user*, int **part*, const int & *count*);

Zoltan_DD_Update takes a list of GIDs and corresponding lists of optional local IDs, optional user data, and optional parts. This routine updates the information for existing directory entries or creates a new entry (filled with given data) if a GID is not found. NULL lists should be passed for optional arguments not desired. This function should be called initially and whenever objects are migrated to keep the distributed directory current.

The user can set the debug level argument in **Zoltan_DD_Create** to determine the module's response to multiple updates for any GID within one update cycle.

Arguments:

dd Distributed directory structure state information.
gid List of GIDs to update (in).
lid List of corresponding local IDs (optional) (in).
user List of corresponding user data (optional) (in).
part List of corresponding parts (optional) (in).
count Number of GIDs in update list.

Returned Value:

int [Error code](#).

C: int **Zoltan_DD_Find** (struct Zoltan_DD_Struct **dd*, [ZOLTAN_ID_PTR](#) *gid*, [ZOLTAN_ID_PTR](#) *lid*, [ZOLTAN_ID_PTR](#) *data*, int **part*, int *count*, int **owner*);
C++: int **Zoltan_DD::Find**([ZOLTAN_ID_PTR](#) *gid*, [ZOLTAN_ID_PTR](#) *lid*, [ZOLTAN_ID_PTR](#) *data*, int **part*, const int & *count*, int **owner*) const;

Given a list of GIDs, **Zoltan_DD_Find** returns corresponding lists of the GIDs' owners, local IDs, parts, data owners, and optional user data. NULL lists must be provided for optional information not being used.

Arguments:

<i>dd</i>	Distributed directory structure state information.
<i>gid</i>	List of GIDs whose information is requested.
<i>lid</i>	Corresponding list of local IDs (optional) (out).
<i>data</i>	Corresponding list of user data (optional) (out).
<i>part</i>	Corresponding list of parts (optional) (out).
<i>count</i>	Count of GIDs in above list.
<i>owner</i>	Corresponding list of data owners (out).

Returned Value:

int	Error code .
-----	------------------------------

ZOLTAN_OK is the normal return.
ZOLTAN_WARN is returned when at least one GID in the *gid* list is not found AND debug level > 0.
ZOLTAN_MEMERR is returned whenever memory can not be allocated.
ZOLTAN_FATAL is returned whenever there is a problem with the input arguments (such as *dd* being NULL) or communications error.

```
C:      int Zoltan_DD_Remove (struct Zoltan_DD_Struct *dd, ZOLTAN\_ID\_PTR gid, int count);
C++:    int Zoltan_DD::Remove( ZOLTAN\_ID\_PTR gid, const int & count);
```

Zoltan_DD_Remove takes a list of GIDs and removes all of their information from the distributed directory.

Arguments:

<i>dd</i>	Distributed directory structure state information.
<i>gid</i>	List of GIDs to eliminate from the directory.
<i>count</i>	Number of GIDs to be removed.

Returned Value:

int	Error code .
-----	------------------------------

```
C:      void Zoltan_DD_Set_Hash_Fn (struct Zoltan_DD_Struct *dd, unsigned int (*hash)
(ZOLTAN\_ID\_PTR, int, unsigned int));
C++:    void Zoltan_DD::Set_Hash_Fn( unsigned int (*hash) (ZOLTAN\_ID\_PTR, int, unsigned int));
```

Enables the user to register a new hash function for the distributed directory. (If this routine is not called, the default hash function [Zoltan_Hash](#) will be used automatically.) This hash function determines which processor maintains the distributed directory entry for a given GID. Inexperienced users do not need this routine.

Experienced users may elect to create their own hash function based on their knowledge of their GID naming scheme. The user's hash function must have calling arguments compatible with [Zoltan_Hash](#). The final argument, *nprocs*, is the number of processors in the communicator passed to [Zoltan_DD_Create](#). Consider that a user has defined a hash function, myhash, as

```
extern int total_num_gid;
unsigned int myhash(ZOLTAN\_ID\_PTR gid, int length, unsigned int nproc)
{
    /* Assuming a processor is more likely to query GIDs that are numerically close to the GIDs it owns, */
    /* this hash function tries to store the gid's directory information near the gid's owning processor's
neighborhood. */
    /* GID length is one ; total_num_gid is a global variable with the total number of GIDs in the application. */

    return ((*gid * nproc) / total_num_gid);
}
```

Then the call to register this hash function is:
Zoltan_DD_Set_Hash(dd, myhash);

Arguments:

- dd* Distributed directory structure state information.
- hash* Name of user's hash function.

Returned Value:

- void NONE
-

```
C: void Zoltan_DD_Stats (struct Zoltan_DD_Struct *dd);
C++: void Zoltan_DD::Stats() const;
```

This routine prints out summary information about the local distributed directory. It includes the hash table length, number of GIDs stored in the local directory, the number of linked lists, and the length of the longest linked list. The debug level (set by an argument to **Zoltan_DD_Create** controls this routine's verbosity.

Arguments:

- dd* Distributed directory structure for state information

Returned Value:

- void NONE
-

```
int Zoltan_DD_Set_Neighbor_Hash_Fn1 (struct Zoltan_DD_Struct *dd, int size);
```

This routine associates the first size GIDs to proc 0, the next size to proc 1, etc. It assumes the GIDs are consecutive numbers. It assumes that GIDs primarily stay near their original owner. The GID length is assumed to be 1. GIDs outside of the range are evenly distributed among the processors via modulo(number of processors). This is a model for the user to develop their own similar routine.

Arguments:

- dd* Distributed directory structure state information.
- size* Number of consecutive GIDs associated with a processor.

Returned Value:

- int [Error code](#).
-

```
int Zoltan_DD_Set_Neighbor_Hash_Fn2 (struct Zoltan_DD_Struct *dd, int *proc, int *low, int *high, int n);
```

This routine allows the user to specify a beginning and ending GID "numbers" per directory processor. It assumes that GIDs primarily stay near their original owner. It requires that the numbers of high, low, & proc entries are all n. It assumes the GID length is 1. It is a model for the user to develop their own similar routine. Users should note the registration of a cleanup routine to free local static memory when the distributed directory is destroyed. GIDs outside the range specified by high and low lists are evenly distributed among the processors via modulo (number of processors).

Arguments:

<i>dd</i>	Distributed directory structure state information.
<i>proc</i>	List of processor ids labeling for corresponding high, low value.
<i>low</i>	List of low GID limits corresponding to proc list.
<i>high</i>	List of high GID limits corresponding to proc list.
<i>n</i>	Number of elements in the above lists. Should be number of processors!

Returned Value:

int	Error code.
-----	-----------------------------

```
C:      int Zoltan_DD_Print (struct Zoltan_DD_Struct *dd);
C++:    int Zoltan_DD::Print () const;
```

This utility displays (to stdout) the entire contents of the distributed directory at one line per GID.

Arguments:

<i>dd</i>	Distributed directory structure state information.
-----------	----------------------------------------------------

Returned Value:

int	Error code.
-----	-----------------------------

User's Notes

Because Zoltan places no restrictions on the content or length of GIDs, hashing does not guarantee a balanced distribution of objects in the distributed directory. Note also, the worst case behavior of a hash table lookup is very bad (essentially becoming a linear search). Fortunately, the average behavior is very good! The user may specify their own hash function via [Zoltan_DD_Set_Hash_Fn](#) to improve performance.

This software module is built on top of the Zoltan Communications functions for efficiency. Improvements to the communications library will automatically benefit the distributed directory.

Examples of Zoltan Usage

Examples for each part of the Zoltan library are provided:

- [General use of Zoltan](#)
- [Load-balancing calling sequence](#)
- [Data migration calling sequences](#)
- [Query functions for a simple application](#)

[[Table of Contents](#) | [Next: General Usage Example](#) | [Previous: Distributed Data Directories](#) | [Privacy and Security](#)]

General Usage Example

An example of general Zoltan usage is included below. This is a C language example. Similar C++ examples may be found in the *examples* directory.

In this example, [Zoltan_Initialize](#) is called using the *argc* and *argv* arguments to the main program. Then a pointer to a Zoltan structure is returned by the call to [Zoltan_Create](#). In this example, all processors will be used by Zoltan, as `MPI_COMM_WORLD` is passed to [Zoltan_Create](#) as the communicator.

Several application query functions are then registered with Zoltan through calls to [Zoltan_Set_Fn](#). Parameters are set through calls to [Zoltan_Set_Param](#). The application then performs in computations, including making calls to Zoltan functions and utilities.

Before its execution ends, the application frees memory used by Zoltan by calling [Zoltan_Destroy](#).

```
/* Initialize the Zoltan library */
struct Zoltan_Struct *zz;
float version;
...
Zoltan\_Initialize(argc, argv, &version);
zz = Zoltan\_Create(MPI_COMM_WORLD);

/* Register query functions. */
Zoltan\_Set\_Fn(zz, ZOLTAN\_NUM\_GEOM\_FN\_TYPE,
               (void (*)()) user_return_dimension, NULL);
Zoltan\_Set\_Fn(zz, ZOLTAN\_GEOM\_FN\_TYPE,
               (void (*)()) user_return_coords, NULL);
Zoltan\_Set\_Fn(zz, ZOLTAN\_NUM\_OBJ\_FN\_TYPE,
               (void (*)()) user_return_num_node, NULL);
Zoltan\_Set\_Fn(zz, ZOLTAN\_OBJ\_LIST\_FN\_TYPE,
               (void (*)()) user_return_owned_nodes, NULL);

/* Set some Zoltan parameters. */
Zoltan\_Set\_Param(zz, "debug_level", "4");

/* Perform application computations, call Zoltan, etc. */
...

/* Free Zoltan data structure before ending application. */
Zoltan\_Destroy (&zz);
```

Typical calling sequence for general usage of the Zoltan library.

```
! Initialize the Zoltan library
type(Zoltan_Struct), pointer :: zz
real(Zoltan_FLOAT) version
integer(Zoltan_INT) ierr
...
ierr = Zoltan\_Initialize(version) ! without argc and argv
zz => Zoltan\_Create(MPI_COMM_WORLD)

! Register load-balancing query functions.
! omit data = C NULL
```



```
ierr = Zoltan\_Set\_Fn(zz, ZOLTAN\_NUM\_GEOM\_FN\_TYPE, user_return_dimension)
ierr = Zoltan\_Set\_Fn(zz, ZOLTAN\_GEOM\_FN\_TYPE, user_return_coords)
ierr = Zoltan\_Set\_Fn(zz, ZOLTAN\_NUM\_OBJ\_FN\_TYPE, user_return_num_node)
ierr = Zoltan\_Set\_Fn(zz, ZOLTAN\_OBJ\_LIST\_FN\_TYPE, user_return_owned_nodes)

! Set some Zoltan parameters.
ierr = Zoltan\_Set\_Param(zz, "debug_level", "4")

! Perform application computations, call Zoltan, etc.
...

! Free Zoltan data structure before ending application.
call Zoltan\_Destroy(zz)
```

Fortran version of general usage example.

Load-Balancing Example

An example of the typical calling sequence for load balancing using Zoltan in a finite element application is shown in the [figure](#) below. An application first selects a load-balancing algorithm by setting the [LB METHOD](#) parameter with [Zoltan_Set_Param](#). Next, other parameter values are set by calls to [Zoltan_Set_Param](#). After some computation, load balancing is invoked by calling [Zoltan_LB_Partition](#). The results of the load balancing include the number of nodes to be imported and exported to the processor, lists of global and local IDs of the imported and exported nodes, and source and destination processors of the imported and exported nodes. A returned argument of [Zoltan_LB_Partition](#) is tested to see whether the new decomposition differs from the old one. If the decompositions differ, some sort of data migration is needed to establish the new decomposition; the details of migration are not shown in this [figure](#) but will be addressed in the [migration examples](#). After the data migration is completed, the arrays of information about imported and exported nodes returned by [Zoltan_LB_Partition](#) are freed by a call to [Zoltan_LB_Free_Part](#).

```

char *lb_method;
int new, num_imp, num_exp, *imp_procs, *exp_procs;
int *imp_to_part, *exp_to_part;
int num_gid_entries, num_lid_entries;
ZOLTAN\_ID\_PTR imp_global_ids, exp_global_ids;
ZOLTAN\_ID\_PTR imp_local_ids, exp_local_ids;

/* Set load-balancing method. */
read_load_balancing_info_from_input_file(&lb_method);
Zoltan\_Set\_Param(zz, "LB_METHOD", lb_method);

/* Reset some load-balancing parameters. */
Zoltan\_Set\_Param(zz, "RCB_Reuse", "TRUE");

/* Perform computations */
...
/* Perform load balancing */
Zoltan\_LB\_Partition(zz,&new,&num_gid_entries,&num_lid_entries,
    &num_imp,&imp_global_ids,&imp_local_ids,&imp_procs,&imp_to_part,
    &num_exp,&exp_global_ids,&exp_local_ids,&exp_procs,&exp_to_part);
if (new)
    perform_data_migration(...);

/* Free memory allocated for load-balancing results by Zoltan library */
Zoltan\_LB\_Free\_Part(&imp_global_ids, &imp_local_ids, &imp_procs, &imp_to_part);
Zoltan\_LB\_Free\_Part(&exp_global_ids, &exp_local_ids, &exp_procs, &exp_to_part);
...

```

Typical calling sequence for performing load balancing with the Zoltan library.

```

character(len=3) lb_method
logical new
integer(Zoltan_INT) num_imp, num_exp
integer(Zoltan_INT) num_gid_entries, num_lid_entries
integer(Zoltan_INT), pointer :: imp_procs(:), exp_procs(:)
integer(Zoltan_INT), pointer :: imp_global_ids(:), exp_global_ids(:) ! global IDs
integer(Zoltan_INT), pointer :: imp_local_ids(:), exp_local_ids(:) ! local IDs
integer(Zoltan_INT) ierr

```

```
! Set load-balancing method.
lb_method = "RCB"
ierr = Zoltan\_Set\_Param(zz, "LB_METHOD", lb_method)

! Reset some load-balancing parameters
ierr = Zoltan\_Set\_Param(zz, "RCB_Reuse", "TRUE")

! Perform computations
...
! Perform load balancing
ierr = Zoltan\_LB\_Partition(zz,new,num_gid_entries,num_lid_entries, &
    num_imp,imp_global_ids,imp_local_ids, &
    imp_procs,imp_to_part, &
    num_exp,exp_global_ids,exp_local_ids, &
    exp_procs,exp_to_part)
if (new) then
    perform_data_migration(...)
endif

! Free memory allocated for load-balancing results by Zoltan library
ierr = Zoltan\_LB\_Free\_Part(imp_global_ids, imp_local_ids, imp_procs, imp_to_part);
ierr = Zoltan\_LB\_Free\_Part(exp_global_ids, exp_local_ids, exp_procs, exp_to_part);
...
```

Fortran version of the load-balancing example.

Migration Examples

Data migration using Zoltan's migration tools can be accomplished in two different ways:

[auto-migration](#), or
[user-guided migration](#).

The choice of migration method depends upon the complexity of the application's data. For some applications, only the objects used in balancing must be migrated; no auxiliary data structures must be moved. Particle simulations are examples of such applications; load balancing is based on the number of particles per processor, and only the particles and their data must be moved to establish the new decomposition. For such applications, Zoltan's auto-migration tools can be used. Other applications, such as finite element methods, perform load balancing on, say, the nodes of the finite element mesh, but nodes that are moved to new processors also need to have their connected elements moved to the new processors, and migrated elements may also need "ghost" nodes (i.e., copies of nodes assigned to other processors) to satisfy their connectivity requirements on the new processor. This complex data migration requires a more user-controlled approach to data migration than the auto-migration capabilities Zoltan can provide.

Auto-Migration Example

In the [figure](#) below, an example of the load-balancing calling sequence for a particle simulation using Zoltan's auto-migration tools is shown. The application requests auto-migration by turning on the **AUTO_MIGRATE** option through a call to [Zoltan_Set_Param](#) and registers functions to pack and unpack a particle's data. During the call to [Zoltan_LB_Partition](#), Zoltan computes the new decomposition and, using calls to the packing and unpacking query functions, automatically migrates particles to their new processors. The application then frees the arrays returned by [Zoltan_LB_Partition](#) and can continue computation without having to perform any additional operations for data migration.

```
/* Tell Zoltan to automatically migrate data for the application. */
Zoltan_Set_Param(zz, "AUTO_MIGRATE", "TRUE");

/* Register additional functions for packing and unpacking data */
/* by migration tools. */
Zoltan_Set_Fn(zz, ZOLTAN_OBJ_SIZE_FN_TYPE,
              (void (*)()) user_return_particle_data_size, NULL);
Zoltan_Set_Fn(zz, ZOLTAN_PACK_OBJ_FN_TYPE,
              (void (*)()) user_pack_particle_data, NULL);
Zoltan_Set_Fn(zz, ZOLTAN_UNPACK_OBJ_FN_TYPE,
              (void (*)()) user_unpack_particle_data, NULL);
...
/* Perform computations */
...
/* Perform load balancing AND automatic data migration! */
Zoltan_LB_Partition(zz,&new,&num_gid_entries,&num_lid_entries,
                   &num_imp,&imp_global_ids,&imp_local_ids,&imp_procs,&imp_to_part,
                   &num_exp,&exp_global_ids,&exp_local_ids,&exp_procs,&exp_to_part);

/* Free memory allocated for load-balancing results by Zoltan */
Zoltan_LB_Free_Part(&imp_global_ids, &imp_local_ids, &imp_procs, &imp_to_part);
```

```
Zoltan\_LB\_Free\_Part(&exp_global_ids, &exp_local_ids, &exp_procs, &exp_to_part);
...
```

Typical calling sequence for using the migration tools' auto-migration capability with the dynamic load-balancing tools.

User-Guided Migration Example

In the following [figure](#), an example of user-guided migration using Zoltan's migration tools for a finite element application is shown. Several migration steps are needed to completely rebuild the application's data structures for the new decomposition. On each processor, newly imported nodes need copies of elements containing those nodes. Newly imported elements, then, need copies of "ghost" nodes, nodes that are in the element but are assigned to other processors. Each of these entities (nodes, elements, and ghost nodes) can be migrated in separate migration steps using the functions provided in the migration tools. First, the assignment of nodes to processors returned by [Zoltan_LB_Partition](#) is established. Query functions that pack and unpack nodes are registered and [Zoltan_Migrate](#) is called using the nodal decomposition returned from [Zoltan_LB_Partition](#). [Zoltan_Migrate](#) packs the nodes to be exported, sends them to other processors, and unpacks nodes received by a processor. The packing routine *migrate_node_pack* includes with each node a list of the element IDs for elements containing that node. The unpacking routine *migrate_node_unpack* examines the list of element IDs and builds a list of requests for elements the processor needs but does not already store. At the end of the nodal migration, each processor has a list of element IDs for elements that it needs to support imported nodes but does not already store. Through a call to [Zoltan_Invert_Lists](#), each processor computes the list of elements it has to send to other processors to satisfy their element requests. Packing and unpacking routines for elements are registered, and [Zoltan_Migrate](#) is again used to move element data to new processors. Requests for ghost nodes can be built within the element packing and unpacking routines, and calls to [Zoltan_Invert_Lists](#) and [Zoltan_Migrate](#), with node packing and unpacking, satisfy requests for ghost nodes. In all three phases of migration, the migration tools handle communication; the application is responsible only for packing and unpacking data and for building the appropriate request lists.

```
/* Assume Zoltan returns a decomposition of the */
/* nodes of a finite element mesh. */
Zoltan\_LB\_Partition(zz,&new,&num_gid_entries,&num_lid_entries,
    &num_imp,&imp_global_ids,&imp_local_ids,&imp_procs,&imp_to_part,
    &num_exp,&exp_global_ids,&exp_local_ids,&exp_procs,&exp_to_part);

/* Migrate the nodes as directed by the results of Zoltan_LB_Partition. */
/* While unpacking nodes, build list of requests for elements needed */
/* to support the imported nodes.*/
Zoltan\_Set\_Fn(zz, ZOLTAN\_OBJ\_SIZE\_FN\_TYPE,
    (void (*)()) migrate_node_size, NULL);
Zoltan\_Set\_Fn(zz, ZOLTAN\_PACK\_OBJ\_FN\_TYPE,
    (void (*)()) migrate_pack_node, NULL);
Zoltan\_Set\_Fn(zz, ZOLTAN\_UNPACK\_OBJ\_FN\_TYPE,
    (void (*)()) migrate_unpack_node, NULL);
Zoltan\_Migrate(zz,num_import,imp_global_ids,imp_local_ids,imp_procs,imp_to_part,
    num_export,exp_global_ids,exp_local_ids,exp_procs,exp_to_part);

/* Prepare for migration of requested elements. */
Zoltan\_Set\_Fn(zz, ZOLTAN\_PACK\_OBJ\_FN\_TYPE,
    (void (*)()) migrate_pack_element, NULL);
Zoltan\_Set\_Fn(zz, ZOLTAN\_UNPACK\_OBJ\_FN\_TYPE,
    (void (*)()) migrate_unpack_element, NULL);
Zoltan\_Set\_Fn(zz, ZOLTAN\_OBJ\_SIZE\_FN\_TYPE,
    (void (*)()) migrate_element_size, NULL);
```

```
/* From the request lists, a processor knows which elements it needs */
/* to support the imported nodes; it must compute which elements to */
/* send to other processors. */
Zoltan_Invert_Lists(zz, Num_Elt_Requests, Elt_Requests_Global_IDs,
    Elt_Requests_Local_IDs, Elt_Requests_Procs, Elt_Requests_to_Part,
    &num_tmp_exp, &tmp_exp_global_ids,
    &tmp_exp_local_ids, &tmp_exp_procs, &tmp_exp_to_part);

/* Processor now knows which elements to send to other processors. */
/* Send the requested elements. While unpacking elements, build */
/* request lists for "ghost" nodes needed by the imported elements. */
Zoltan_Migrate(zz, Num_Elt_Requests, Elt_Requests_Global_IDs,
    Elt_Requests_Local_IDs, Elt_Requests_Procs, Elt_Request_to_Part,
    num_tmp_exp_objs, tmp_exp_global_ids,
    tmp_exp_local_ids, tmp_exp_procs, tmp_exp_to_part);

/* Repeat process for "ghost" nodes. */
...
```

Typical calling sequence for user-guided use of the migration tools in Zoltan.

Query-Function Examples

Examples of query functions provided by a simple application are included below. The general-interface examples include a simple implementation of [ZOLTAN_GEOM_FN](#) and [ZOLTAN_OBJ_LIST_FN](#) query functions and variants of the simple implementation that exploit local identifiers and data pointers. Migration examples for packing and unpacking objects are also included. Robust error checking is not included in the routines; application developers should include more explicit error checking in their query functions.

- [General Interface Examples](#)
 - [Basic example](#)
 - [User-defined data pointer](#)
- [Migration Examples](#)
 - [Packing and unpacking functions](#)

All the examples use a mesh data structure consisting of nodes in the mesh. these nodes are the objects passed to Zoltan. A node is described by its 3D coordinates and a global ID number that is unique across all processors. The type definitions for the mesh and node data structures used in the examples are included [below](#).

```

/* Node data structure. */
/* A node consists of its 3D coordinates and */
/* an ID number that is unique across all processors. */
struct Node_Type {
    double Coordinates[3];
    int Global_ID_Num;
};

/* Mesh data structure. */
/* Mesh consists of an array of nodes and */
/* the number of nodes owned by the processor. */
struct Mesh_Type {
    struct Node_Type Nodes[MAX_NODES];
    int Number_Owned;
};
    
```

Data types for the query-function examples.

```

! Node data structure.
! A node consists of its 3D coordinates and
! an ID number that is unique across all processors.
type Node_Type
    real(Zoltan_DOUBLE) :: Coordinates(3)
    integer(Zoltan_INT) :: Global_ID_Num
end type Node_Type

! Mesh data structure.
! Mesh consists of an array of nodes and
! the number of nodes owned by the processor.
type Mesh_Type
    type(Node_Type) :: Nodes(MAX_NODES)
    integer(Zoltan_INT) :: Number_Owned
    
```



```
end type Mesh_Type
```

Data types for the Fortran query-function examples.

General Interface Query Function Examples

In the following examples, [ZOLTAN_OBJ_LIST_FN](#) and [ZOLTAN_GEOM_FN](#) query functions are implemented for an application using the mesh and node data structures described [above](#). The nodes are the objects passed to Zoltan.

Through a call to [Zoltan_Set_Fn](#), the function *user_return_owned_nodes* is registered as the [ZOLTAN_OBJ_LIST_FN](#) query function. It returns global and local identifiers for each node owned by a processor.

The function *user_return_coords* is registered as a [ZOLTAN_GEOM_FN](#) query function. Given the global and local identifiers for a node, this function returns the node's coordinates. All the examples exploit the local identifier to quickly locate nodal data. If such an identifier is not available in an application, a search using the global identifier can be performed.

The [Basic Example](#) includes the simplest implementation of the query routines. In the query routines, it uses global application data structures and a local numbering scheme for the local identifiers. The [User-Defined Data Pointer Example](#) uses only local application data structures; this model is useful if the application does not have global data structures or if objects from more than one data structure are to be passed to Zoltan. Differences between the latter example and the Basic Example are shown in [red](#).

Basic Example

In the simplest example, the query functions access the application data through a global data structure (*Mesh*) representing the mesh. In the calls to [Zoltan_Set_Fn](#), no pointers to application data are registered with the query function (i.e., the *data* pointer is not used). A node's local identifier is an integer representing the index in the *Mesh.Nodes* array of the node. The local identifier is set to the index's value in *user_return_owned_nodes*. It is used to access the global *Mesh.Nodes* array in *user_return_coords*.

```
/* in application's program file */
#include "zoltan.h"

/* Declare a global Mesh data structure. */
struct Mesh_Type Mesh;

main()
{
...
    /* Indicate that local and global IDs are one integer each. */
    Zoltan_Set_Param(zz, "NUM_GID_ENTRIES", "1");
    Zoltan_Set_Param(zz, "NUM_LID_ENTRIES", "1");

    /* Register query functions. */
    /* Do not register a data pointer with the functions; */
    /* the global Mesh data structure will be used. */
    Zoltan_Set_Fn(zz, ZOLTAN_GEOM_FN_TYPE,
        (void (*)()) user_return_coords, NULL);
    Zoltan_Set_Fn(zz, ZOLTAN_OBJ_LIST_FN_TYPE,
        (void (*)()) user_return_owned_nodes, NULL);
...
}
```

```

}

void user_return_owned_nodes(void *data,
    int num_gid_entries, int num_lid_entries,
    ZOLTAN\_ID\_PTR global_ids, ZOLTAN\_ID\_PTR local_ids,
    int wgt_dim, float *obj_wgts,
    int *ierr)
{
    int i;
    /* return global node numbers as global_ids. */
    /* return index into Nodes array for local_ids. */
    for (i = 0; i < Mesh.Number_Owned; i++){
        global_ids[i*num_gid_entries] = Mesh.Nodes[i].Global_ID_Num;
        local_ids[i*num_lid_entries] = i;
    }
    *ierr = ZOLTAN_OK;
}

void user_return_coords(void *data,
    int num_gid_entries, int num_lid_entries,
    ZOLTAN\_ID\_PTR global_id, ZOLTAN\_ID\_PTR local_id,
    double *geom_vec, int *ierr)
{
    /* use local_id to index into the Nodes array. */
    geom_vec[0] = Mesh.Nodes[local_id[0]].Coordinates[0];
    geom_vec[1] = Mesh.Nodes[local_id[0]].Coordinates[1];
    geom_vec[2] = Mesh.Nodes[local_id[0]].Coordinates[2];
    *ierr = ZOLTAN_OK;
}

```

Example of general interface query functions (simplest implementation).

```

! in application's program file

module Global_Mesh_Data
! Declare a global Mesh data structure.
    type(Mesh_Type) :: Mesh
end module

program query_example_1
use zoltan
...
    ! Indicate that local and global IDs are one integer each.
    ierr = Zoltan\_Set\_Param(zz, "NUM\_GID\_ENTRIES", "1");
    ierr = Zoltan\_Set\_Param(zz, "NUM\_LID\_ENTRIES", "1");

    ! Register query functions.
    ! Do not register a data pointer with the functions;
    ! the global Mesh data structure will be used.
    ierr = Zoltan\_Set\_Fn(zz, ZOLTAN\_GEOM\_FN\_TYPE, user_return_coords)
    ierr = Zoltan\_Set\_Fn(zz, ZOLTAN\_OBJ\_LIST\_FN\_TYPE, user_return_owned_nodes)
...
end program

subroutine user_return_owned_nodes(data, &
    num_gid_entries, num_lid_entries, &

```

```

    global_ids, local_ids, wgt_dim, obj_wgts, ierr)
use zoltan
use Global_Mesh_Data
integer(Zoltan_INT) :: data(1) ! dummy declaration, do not use
integer(Zoltan_INT), intent(in) :: num_gid_entries, num_lid_entries
integer(Zoltan_INT), intent(out) :: global_ids(*), local_ids(*)
integer(Zoltan_INT), intent(in) :: wgt_dim
real(Zoltan_FLOAT), intent(out) :: obj_wgts(*)
integer(Zoltan_INT), intent(out) :: ierr
integer i
    ! return global node numbers as global_ids.
    ! return index into Nodes array for local_ids.
do i = 1, Mesh%Number_Owned
    global_ids(1+(i-1)*num_gid_entries) = &
        Mesh%Nodes(i)%Global_ID_Num
    local_ids(1+(i-1)*num_lid_entries) = i
end do
ierr = ZOLTAN_OK
end subroutine

subroutine user_return_coords(data, num_gid_entries, num_lid_entries, &
    global_id, local_id, geom_vec, ierr)
use zoltan
use Global_Mesh_Data
integer(Zoltan_INT) :: data(1) ! dummy declaration, do not use
integer(Zoltan_INT), intent(in) :: num_gid_entries, num_lid_entries
integer(Zoltan_INT), intent(in) :: global_id(*), local_id(*)
real(Zoltan_DOUBLE), intent(out) :: geom_vec(*)
integer(Zoltan_INT), intent(out) :: ierr
    ! use local_id to index into the Nodes array.
    geom_vec(1:3) = Mesh%Nodes(local_id(1))%Coordinates
    ierr = ZOLTAN_OK
end subroutine

```

Fortran example of general interface query functions (simplest implementation).

User-Defined Data Pointer Example

In this example, the address of a local mesh data structure is registered with the query functions for use by those functions. This change eliminates the need for a global mesh data structure in the application. The address of the local data structure is included as an argument in calls to [Zoltan_Set_Fn](#). This address is then used in *user_return_owned_nodes* and *user_return_coords* to provide data for these routines. It is cast to the [Mesh_Type](#) data type and accessed with local identifiers as in the [Basic Example](#). Differences between this example and the [Basic Example](#) are shown in **red**.

This model is useful when the application does not have a global data structure that can be accessed by the query functions. It can also be used for operations on different data structures. For example, if an application had more than one mesh, load balancing could be performed separately on each mesh without having different query routines for each mesh. Calls to [Zoltan_Set_Fn](#) would define which mesh should be balanced, and the query routines would access the mesh currently designated by the [Zoltan_Set_Fn](#) calls.

```

/* in application's program file */
#include "zoltan.h"

```

```

main()
{
    /* declare a local mesh data structure. */
    struct Mesh_Type mesh;

    ...
    /* Indicate that local and global IDs are one integer each. */
    Zoltan_Set_Param(zz, "NUM_GID_ENTRIES", "1");
    Zoltan_Set_Param(zz, "NUM_LID_ENTRIES", "1");

    /* Register query functions. */
    /* Register the address of mesh as the data pointer. */
    Zoltan_Set_Fn(zz, ZOLTAN_GEOM_FN_TYPE,
        (void (*)()) user_return_coords, &mesh);
    Zoltan_Set_Fn(zz, ZOLTAN_OBJ_LIST_FN_TYPE,
        (void (*)()) user_return_owned_nodes, &mesh);
    ...
}

void user_return_owned_nodes(void *data,
    int num_gid_entries, int num_lid_entries,
    ZOLTAN_ID_PTR global_ids, ZOLTAN_ID_PTR local_ids,
    int wgt_dim, float *obj_wgts,
    int *ierr)
{
    int i;
    /* cast data pointer to type Mesh_Type. */
    struct Mesh_Type *ptr = (struct Mesh_Type *) data;

    /* return global node numbers as global_ids. */
    /* return index into Nodes array for local_ids. */
    for (i = 0; i < ptr->Number_Owned; i++) {
        global_ids[i*num_gid_entries] = ptr->Nodes[i].Global_ID_Num;
        local_ids[i*num_lid_entries] = i;
    }
    *ierr = ZOLTAN_OK;
}

void user_return_coords(void *data,
    int num_gid_entries, int num_lid_entries,
    ZOLTAN_ID_PTR global_id, ZOLTAN_ID_PTR local_id,
    double *geom_vec, int *ierr)
{
    /* cast data pointer to type Mesh_Type. */
    struct Mesh_Type *ptr = (struct Mesh_Type *) data;

    /* use local_id to address the requested node. */
    geom_vec[0] = ptr->Nodes[local_id[0]].Coordinates[0];
    geom_vec[1] = ptr->Nodes[local_id[0]].Coordinates[1];
    geom_vec[2] = ptr->Nodes[local_id[0]].Coordinates[2];
    *ierr = ZOLTAN_OK;
}

```

Example of general interface query functions using the application-defined data pointer.

```

/* included in file zoltan_user_data.f90 */

```

```

! User defined data type as wrapper for Mesh
type Zoltan_User_Data_1
    type(Mesh_Type), pointer :: ptr
end type Zoltan_User_Data_1

! in application's program file

program query_example_3
use zoltan
! declare a local mesh data structure and a User_Data to point to it.
type(Mesh_Type), target :: mesh
type(Zoltan_User_Data_1) data
...
    ! Indicate that local and global IDs are one integer each.
    ierr = Zoltan_Set_Param(zz, "NUM_GID_ENTRIES", "1");
    ierr = Zoltan_Set_Param(zz, "NUM_LID_ENTRIES", "1");

    ! Register query functions.
    ! Use the User_Data variable to pass the mesh data
    data%ptr => mesh
    ierr = Zoltan_Set_Fn(zz, ZOLTAN_GEOM_FN_TYPE, user_return_coords, data)
    ierr = Zoltan_Set_Fn(zz, ZOLTAN_OBJ_LIST_FN_TYPE,
        user_return_owned_nodes, data)
...
end program

subroutine user_return_owned_nodes(data, &
    num_gid_entries, num_lid_entries, &
    global_ids, local_ids, wgt_dim, obj_wgts, ierr)
use zoltan
type(Zoltan_User_Data_1) :: data
integer(Zoltan_INT), intent(in) :: num_gid_entries, num_lid_entries
integer(Zoltan_INT), intent(out) :: global_ids(*), local_ids(*)
integer(Zoltan_INT), intent(in) :: wgt_dim
real(Zoltan_FLOAT), intent(out) :: obj_wgts(*)
integer(Zoltan_INT), intent(out) :: ierr
integer i
type(Mesh_Type), pointer :: Mesh

! extract the mesh from the User_Data argument
Mesh => data%ptr

! return global node numbers as global_ids.
! return index into Nodes array for local_ids.
do i = 1, Mesh%Number_Owned
    global_ids(1+(i-1)*num_gid_entries) = &
        Mesh%Nodes(i)%Global_ID_Num
    local_ids(1+(i-1)*num_lid_entries) = i
end do
ierr = ZOLTAN_OK
end subroutine

subroutine user_return_coords(data, global_id, local_id, &
    geom_vec, ierr)
use zoltan
type(Zoltan_User_Data_1) :: data
integer(Zoltan_INT), intent(in) :: num_gid_entries, num_lid_entries

```

```
integer(Zoltan_INT), intent(in) :: global_id(*), local_id(*)
real(Zoltan_DOUBLE), intent(out) :: geom_vec(*)
integer(Zoltan_INT), intent(out) :: ierr
type(Mesh_Type), pointer :: Mesh

! extract the mesh from the User_Data argument
Mesh => data%ptr

! use local_id to index into the Nodes array.
geom_vec(1:3) = Mesh%Nodes(local_id(1))%Coordinates
ierr = ZOLTAN_OK
end subroutine
```

Fortran example of general interface query functions using the application-defined data pointer.

Migration Examples

Packing and Unpacking Data

Simple migration query functions for the [Basic Example](#) are included [below](#). These functions are used by the migration tools to move nodes among the processors. The functions `user_size_node`, `user_pack_node`, and `user_unpack_node` are registered through calls to [Zoltan_Set_Fn](#). Query function `user_size_node` returns the size (in bytes) of data representing a single node. Query function `user_pack_node` copies a given node's data into the communication buffer `buf`. Query function `user_unpack_node` copies a data for one node from the communication buffer `buf` into the `Mesh.Nodes` array on its new processor.

These query routines are simple because the application does not dynamically allocate memory for each node. Such dynamic allocation would have to be accounted for in the [ZOLTAN_OBJ_SIZE_FN](#), [ZOLTAN_PACK_OBJ_FN](#), and [ZOLTAN_UNPACK_OBJ_FN](#) routines.

```
main()
{
...
/* Register migration query functions. */
/* Do not register a data pointer with the functions; */
/* the global Mesh data structure will be used. */
Zoltan_Set_Fn(zz, ZOLTAN_OBJ_SIZE_FN_TYPE,
              (void (*)()) user_size_node, NULL);
Zoltan_Set_Fn(zz, ZOLTAN_PACK_OBJ_FN_TYPE,
              (void (*)()) user_pack_node, NULL);
Zoltan_Set_Fn(zz, ZOLTAN_UNPACK_OBJ_FN_TYPE,
              (void (*)()) user_unpack_node, NULL);
...
}

int user_size_node(void *data,
                  int num_gid_entries, int num_lid_entries,
                  ZOLTAN_ID_PTR global_id, ZOLTAN_ID_PTR local_id, int *ierr)
{
/* Return the size of data associated with one node. */
/* This case is simple because all nodes have the same size. */
*ierr = ZOLTAN_OK;
return(sizeof(struct Node_Type));
}
```

```
void user_pack_node(void *data,
    int num_gid_entries, int num_lid_entries,
    ZOLTAN\_ID\_PTR global_id, ZOLTAN\_ID\_PTR local_id,
    int dest_proc, int size, char *buf, int *ierr)
{
    /* Copy the specified node's data into buffer buf. */
    struct Node_Type *node_buf = (struct Node_Type *) buf;

    *ierr = ZOLTAN_OK;
    node_buf->Coordinates[0] = Mesh.Nodes[local_id[0]].Coordinates[0];
    node_buf->Coordinates[1] = Mesh.Nodes[local_id[0]].Coordinates[1];
    node_buf->Coordinates[2] = Mesh.Nodes[local_id[0]].Coordinates[2];
    node_buf->Global_ID_Num = Mesh.Nodes[local_id[0]].Global_ID_Num;
}

void user_unpack_node(void *data, int num_gid_entries,
    ZOLTAN\_ID\_PTR global_id, int size,
    char *buf, int *ierr)
{
    /* Copy the node data in buf into the Mesh data structure. */
    int i;
    struct Node_Type *node_buf = (struct Node_Type *) buf;

    *ierr = ZOLTAN_OK;
    i = Mesh.Number_Owned;
    Mesh.Number_Owned = Mesh.Number_Owned + 1;
    Mesh.Nodes[i].Coordinates[0] = node_buf->Coordinates[0];
    Mesh.Nodes[i].Coordinates[1] = node_buf->Coordinates[1];
    Mesh.Nodes[i].Coordinates[2] = node_buf->Coordinates[2];
    Mesh.Nodes[i].Global_ID_Num = node_buf->Global_ID_Num;
}
```

Example of migration query functions for the [Basic Example](#).

Release Notes

Release notes are available for the following releases of Zoltan:

[Zoltan Release v3.3](#)
[Zoltan Release v3.2](#)
[Zoltan Release v3.1](#)
[Zoltan Release v3.0](#)
[Zoltan Release v2.1](#)
[Zoltan Release v2.02](#)
[Zoltan Release v2.01](#)
[Zoltan Release v2.0](#)
[Zoltan Release v1.54](#)
[Zoltan Release v1.53](#)
[Zoltan Release v1.52](#)
[Zoltan Release v1.5](#)
[Zoltan Release v1.3](#)

Zoltan Release Notes v3.3: July 31, 2010

Features:

- New [local ordering method based on space-filling curves](#) to improve memory and cache locality within a processor.
 - Ability to call graph partitioning algorithms using hypergraph callback functions; this capability is useful applications with, say, block-structured matrix distributions (e.g., SuperLU), where all information about a matrix row or column is not available on a single processor.
 - Improved execution time of parallel hypergraph partitioning.
-

Zoltan Release Notes v3.2: September 24, 2009

Features:

- New [interface](#) to [Scotch](#) and [PT-Scotch](#) parallel graph partitioning algorithms.
- Simplified interface to graph [ordering](#) and [coloring](#) algorithms
- Automated symmetrization of graphs for graph partitioning, coloring and ordering. (See parameters GRAPH_SYMMETRIZE and GRAPH_SYM_WEIGHT in the [Scotch](#) and [ParMETIS](#) graph packages.)
- Improved function [Zoltan_LB_Eval](#) returns more information about a decomposition to users.
- Improved examples showing Zoltan usage in C and C++ are included in *zoltan/example*.
- Improved support for [builds under autotools](#), including builds of Zoltan's F90 interface.
- New support for [CMake builds](#) and testing through Trilinos; builds of Zoltan's F90 interface are included.
- Improved integration into [Isorropia](#) partitioners for Trilinos' Epetra classes.

Backward compatibility:

- Interfaces to [Zoltan_Color](#), [Zoltan_Order](#) and [Zoltan_LB_Eval](#) have changed.
- The Zoltan native build environment, while still distributed, will no longer be supported. Users should use the [autotools or CMake](#) systems. Builds of the Zoltan F90 interface are supported in both autotools and CMake.

Zoltan Release Notes v3.1: September 26, 2008

Zoltan v3.1 includes the following new features:

- Important new capabilities for Matrix ordering are included in Zoltan v3.1.
 - A graph/matrix ordering interface to [PT-Scotch](#), a high-quality graph ordering and partitioning library from the University of Bordeaux, is now available.
 - Zoltan's new [matrix ordering interface](#) returns ordering information such as permutations and separators to the application.
- New [hypergraph partitioning](#) options for inexpensively [refining](#) partitions are included and controlled simply by parameters [LB_APPROACH](#) and [PHG_MULTILEVEL](#).
- Robustness improvements have been made to Zoltan's parallel [hypergraph](#) partitioner and repartitioner.
- A new [Autotools build environment](#) is available. The native Zoltan build environment is still supported in this release.
- [Serial, non-MPI builds](#) of Zoltan are enabled through the new Autotools build environment.
- Zoltan is now a [Trilinos](#) package. Integration with [Trilinos](#), is now tighter and more seamless. In particular, Zoltan is the default partitioner for the Trilinos package [Isorropia](#), a matrix-based interface to Zoltan.

Please see the [backward compatibility](#) section for a detailed description of changes that may affect current users.

Zoltan Release Notes v3.0: May 1, 2007

Zoltan v3.0 includes major new features.

- [Parallel Hypergraph Repartitioning](#) combining the improved communication metric of [hypergraph partitioning](#) with a [new model](#) for representing an existing partition while computing a new one. This work received the "Best Algorithms Paper Award" at the [2007 IEEE International Parallel and Distributed Processing Symposium](#).
- [Hypergraph refinement](#) to quickly improve the quality of an existing partition.
- Improved partition quality within the Zoltan [parallel hypergraph partitioner](#).
- [Parallel graph partitioning](#) using Zoltan's parallel hypergraph partitioner.
- Hypergraph partitioning with [fixed vertices](#) that allows application to assign or "fix" objects to a desired part before partitioning.
- Improved [part remapping](#) to reduce data migration costs in all partitioners.
- Hybrid [hierarchical partitioning](#) that allows different partitioning algorithms to be applied within a hierarchy of computers (e.g., partitioning across a cluster of shared-memory processors, followed by partitioning within each shared-memory processor).
- A new scheme for more easily specifying partitioning [methods](#) and [approaches](#) within the hypergraph and graph partitioners.
- Very [simple partitioners](#) that serve as testing tools and examples for usage.

Please see the [backward compatibility](#) section for a detailed description of changes that may affect current users.

Zoltan Release Notes v2.1: October 5, 2006

Zoltan v2.1 includes a significant bug fix for the hypergraph partitioner. We strongly recommend that users upgrade to Zoltan v2.1.

Zoltan Release Notes v2.02: September 26, 2006

Zoltan v2.02 includes bugfixes:

- Zoltan_LB_Eval now correctly computes edge cuts with respect to parts when parts are spread across more than one processor.
 - Extraneous (and annoying) print statement has been removed from Zoltan partitioning method RCB.
-

Zoltan Release Notes v2.01: August 2006

Zoltan v2.01 includes enhancements to version 2.0.

- F90 interface fixes to comply with standard F90 (e.g., shortened variable names and continuation lines). The hypergraph callback function names have changed, but C and C++ compatibility with v2.0 is maintained.
 - Performance improvement to initial building of hypergraphs from application data.
 - Major bug fix for dense-edge removal in parallel hypergraph method; partitioning of hypergraphs with edges containing more than 25% of the vertices was affected by this bug.
 - Minor fixes to parallel hypergraph code.
-

Zoltan Release Notes v2.0: April 2006

Zoltan v2.0 includes several major additions:

- [Parallel hypergraph partitioning](#).
 - [Parallel graph coloring](#), both distance-1 and distance-2.
 - [Multicriteria geometric partitioning \(RCB\)](#).
 - [C++ interface](#).
-

Zoltan Release Notes v1.54

Some versions of MPICH have a bug in MPI_Reduce_scatter; they can report errors with MPI_TYPE_INDEXED. In Zoltan v1.54's unstructured communication package, calls to MPI_Reduce_scatter have been replaced with separate calls to MPI_Reduce and MPI_Scatter.

Zoltan Release Notes v1.53

Zoltan v1.53 includes the following new capabilities:

- Portability to BSD Unix and Mac OS X was added.
 - Averaging of RCB and RIB cuts was added; see Zoltan parameter [AVERAGE_CUTS](#).
 - A new function [Zoltan_RCB_Box](#) returns information about subdomain bounding boxes in RCB decompositions.
 - F90 interface to [Zoltan_Order](#) was added.
 - Warnings that load-imbalance tolerance was not met are no longer printed when [DEBUG_LEVEL](#) == 0.
 - Minor bugs were addressed.
-

Zoltan Release Notes v1.52

Zoltan v1.52 includes the following new capabilities:

- List-based graph callback functions [ZOLTAN_NUM_EDGES_MULTI_FN](#) and [ZOLTAN_EDGE_LIST_MULTI_FN](#) were added to mirror support and performance given by the list-based geometric function [ZOLTAN_GEOM_MULTI_FN](#).
- Support for ParMETIS v3.1 was added.
- Minor bugs were addressed.

Zoltan Release Notes v1.5

This section describes improvements to Zoltan in Version 1.5. Every attempt was made to keep Zoltan v1.3 backwardly compatible with previous versions. Users of previous versions of Zoltan should refer to the [Backward Compatibility Notes](#).

Short descriptions of the following features are included below; follow the links for more details.

- [Part remapping](#)
- [Unequal Numbers of Parts and Processors](#)
- [Non-Uniform Part Sizes](#)
- [Zoltan Interface Updated](#)
- [Robust HSFC Box Assign](#)
- [Matrix Ordering](#)
- [Performance Improvements](#)
- [Bug Fixes](#)

Part Remapping

During partitioning, Zoltan v1.5 can renumber parts so that the input and output partitions have greater overlap (and, thus, lower data-migration costs). This remapping is controlled by Zoltan parameter [REMAP](#). Experiments have shown that using this parameter can greatly reduce data migration costs.

Unequal Numbers of Parts and Processors

Zoltan v1.5 can be used to generate k parts on p processors, where k is not equal to p . Function [Zoltan_LB_Partition](#) (replacing [Zoltan_LB_Balance](#)) can generate arbitrary numbers of parts on the given processors. The number of desired parts is set with parameters [NUM_GLOBAL_PARTS](#) or [NUM_LOCAL_PARTS](#). Both part and processor information are returned by [Zoltan_LB_Partition](#), [Zoltan_LB_Box_PP_Assign](#), and [Zoltan_LB_Point_PP_Assign](#). New Zoltan query functions [ZOLTAN_PART_FN](#) and [ZOLTAN_PART_MULTI_FN](#) return objects' part information to Zoltan. [Zoltan_LB_Balance](#) can still be used for k equal to p .

Non-Uniform Part Sizes

Part sizes for local and global parts can be specified using [Zoltan_LB_Set_Part_Sizes](#), allowing non-uniformly sized parts to be generated by Zoltan's partitioning algorithms.

Zoltan Interface Updated

To support the concept of parts separate from processors, many new interface functions were added to Zoltan v1.5 (e.g., [Zoltan_LB_Partition](#) and [Zoltan_Migrate](#)). These functions mimic previous Zoltan functions (e.g., [Zoltan_LB_Balance](#) and [Zoltan_Help_Migrate](#), respectively), but include both part and processor information. Both the new and old interface functions work in Zoltan v1.5. See the notes on [Backward Compatibility](#).

Robust HSFC Box Assign

Function [Zoltan_LB_Box_PP_Assign](#) now works for the [Hilbert Space-Filling Curve algorithm \(HSFC\)](#), in addition to the [RCB](#) and [RIB](#) algorithms supported in previous versions of Zoltan. [Zoltan_LB_Point_PP_Assign](#) continues to work for [HSFC](#), [RCB](#) and [RIB](#).

Matrix Ordering

Zoltan v1.5 contains a matrix-ordering interface [Zoltan_Order](#) to ParMETIS' matrix-ordering functions. New graph-based matrix-ordering algorithms can be easily added behind this interface.

Performance Improvements

Many performance improvements were added to Zoltan v1.5.

- List-based callback functions have been added to Zoltan ([ZOLTAN_GEOM_MULTI_FN](#), [ZOLTAN_PART_MULTI_FN](#), [ZOLTAN_OBJ_SIZE_MULTI_FN](#), [ZOLTAN_PACK_OBJ_MULTI_FN](#), and [ZOLTAN_UNPACK_OBJ_MULTI_FN](#)); these functions allow entire lists of data to be passed from the application to Zoltan, replacing per-object callbacks.
- [Zoltan_Migrate](#) now can accept either import lists, export lists, or both. It is no longer necessary to call [Zoltan_Invert_Lists](#) or [Zoltan_Compute_Destinations](#) to get appropriate input for [Zoltan_Migrate](#).
- Zoltan v1.5 contains performance improvements within individual algorithms. We recommend users upgrade to the latest version.

Bug Fixes

Bug fixes were made to Zoltan's algorithms and interface. Users of previous versions of Zoltan are encouraged to upgrade.

Zoltan Release Notes v1.3

This section describes improvements to Zoltan in Version 1.3. Every attempt was made to keep Zoltan v1.3 backwardly compatible with previous versions. Users of previous versions of Zoltan should refer to the [Backward Compatibility Notes](#).

Short descriptions of the following features are included below; follow the links for more details.

- [More Data Services](#)
- [New Hilbert Space-Filling Curve Partitioning](#)

[Support for Structured-Grid Partitioning](#)
[Support for ParMETIS v3.0](#)
[Performance Improvements](#)
[Zoltan Interface Updated](#)
[Improved Test Suite](#)
[Bug Fixes](#)

More Data Services

Zoltan's mission has been widened beyond its original focus on dynamic load-balancing algorithms. Now Zoltan also provides data management services to parallel, unstructured, and adaptive computations. Several packages of parallel data services have been added and made available to application developers. These services include the following:

- An [unstructured communication package](#) that simplifies complicated communication by insulating applications from the details of message sends and receives.
 - A [distributed data directory](#) that allows applications to efficiently (in memory and time) locate off-processor data.
 - A [dynamic memory management package](#) that simplifies debugging of memory allocation problems on state-of-the-art parallel computers.
-

New Hilbert Space-Filling Curve Partitioning

Zoltan now includes a fast, efficient implementation of [Hilbert Space-Filling Curve \(HSFC\)](#) partitioning. This geometric method also includes support for [Zoltan LB_Box_Assign](#) and [Zoltan LB_Point_Assign](#) functions.

Support for Structured-Grid Partitioning

Zoltan's [Recursive Coordinate Bisection \(RCB\)](#) partitioning algorithm has been enhanced to allow generation of strictly rectilinear subdomains. This capability can be used for partitioning of grids for structured-grid applications. See parameter [RCB_RECTILINEAR_BLOCKS](#).

Support for ParMETIS v3.0

In addition to providing interfaces to [ParMETIS v2.0](#), Zoltan now provides an interfaces [ParMETIS v3.0](#). Full support of ParMETIS v3.0's multiconstraint and multiobjective partitioning is included.

Performance Improvements

Performance of Zoltan's partitioning algorithms has been improved through a number of code optimizations and new features. In addition, user parameter [RETURN_LISTS](#) can be used to specify which returned arguments are computed by [Zoltan LB_Balance](#), allowing reduced work in partitioning. In the [Recursive Coordinate Bisection \(RCB\)](#) partitioning algorithm, user parameters allow cut directions to be locked in an attempt to minimize data movement; see parameters [RCB_LOCK DIRECTIONS](#) and [RCB_SET DIRECTIONS](#).

Zoltan Interface Updated

Zoltan has adopted a more modular design, making it easier to use by applications and easier to modify by algorithm

developers. Names in the [Zoltan interface](#) and code are tied more closely to their functionality. Full [backward compatibility](#) is supported for users of previous versions of Zoltan.

Improved Test Suite

The Zoltan [test suite](#) has been improved, with more tests providing greater code coverage and platform-specific answer files accounting for differences due to computer architectures.

Bug Fixes

Some bug fixes were made to Zoltan's algorithms and interface. Users of previous versions of Zoltan are encouraged to upgrade.

[\[Table of Contents](#) | [Next: Backward Compatibility](#) | [Previous: Query Function Examples](#) | [Privacy and Security\]](#)

Backward Compatibility with Previous Versions of Zoltan

As new features have been added to Zoltan, backward compatibility with previous versions of Zoltan has been maintained. Thus, users of previous versions of Zoltan can upgrade to a new version **without changing their application source code**. Modifications to application source code are needed **only** if the applications use new Zoltan functionality.

Enhancements to the Zoltan interface are described below.

- [Versions 3.2 and higher](#)
- [Versions 3.1 and higher](#)
- [Versions 3.0 and higher](#)
- [Versions 1.5 and higher](#)
- [Versions 1.3 and higher](#)

Backward Compatibility: Versions 3.2 and higher

Interfaces to [Zoltan_Color](#), [Zoltan_Order](#) and [Zoltan_LB_Eval](#) have changed.

The Zoltan native build environment, while still distributed, will no longer be supported. Users should use the [autotools or CMake](#) systems. Builds of the Zoltan F90 interface are supported in both autotools and CMake.

Backward Compatibility: Versions 3.1 and higher

Terminology referring to partitions and parts was clarified. A "partition" describes the entire layout of the data across processes. A "part" is a subset of the data assigned to a single process. A partition is made up of many parts; the set of all the parts is a partition.

We applied this naming convention more consistently throughout Zoltan. Old parameters NUM_GLOBAL_PARTITIONS and NUM_LOCAL_PARTITIONS have been more accurately renamed [NUM_GLOBAL_PARTS](#) and [NUM_LOCAL_PARTS](#). Old query functions ZOLTAN_PARTITIONS_MULTI_FN and ZOLTAN_PARTITION_FN have been more accurately renamed [ZOLTAN_PART_MULTI_FN](#) and [ZOLTAN_PART_FN](#). However, in both cases, the old naming convention still works in the Zoltan library.

Backward Compatibility: Versions 3.0 and higher

A new naming convention was implemented to better categorize partitioning methods. For more details, see parameters

- [LB_METHOD](#),
- [LB_APPROACH](#),
- [GRAPH_PACKAGE](#), and
- [HYPERGRAPH_PACKAGE](#).

Former valid values of LB_METHOD should continue to work. In particular, values of LB_METHOD for geometric partitioners [RCB](#), [RIB](#), [HSFC](#), and [REFTREE](#) are unchanged.

The default graph partitioner has been changed from [ParMETIS](#) to [Zoltan PHG](#). This change was made to provide graph partitioning capability without reliance on the third-party library ParMETIS.

Because Zoltan is designed primarily for dynamic load balancing, The default partitioning approach [LB_APPROACH](#) is now "repartition." This change affects only Zoltan's hypergraph partitioner [PHG](#).

Backward Compatibility: Versions 1.5 and higher

The ability to generate more parts than processors was added to Zoltan in version 1.5. Thus, Zoltan's partitioning and migration routines were enhanced to return and use both part assignments and processor assignments. New interface and query functions were added to support this additional information. All former Zoltan [parameters](#) apply to the new functions as they did to the old; new parameters [NUM_GLOBAL_PARTS](#) and [NUM_LOCAL_PARTS](#) apply only to the new functions.

The table below lists the Zoltan function that uses both part and processor information, along with the analogous function that returns only processor information. Applications requiring only one part per processor can use either version of the functions.

Function with Part and Processor info (v1.5 and higher)	Function with only Processor info (v1.3 and higher)
Zoltan_LB_Partition	Zoltan_LB_Balance
Zoltan_LB_Point_PP_Assign	Zoltan_LB_Point_Assign
Zoltan_LB_Box_PP_Assign	Zoltan_LB_Box_Assign
Zoltan_Invert_Lists	Zoltan_Compute_Destinations
Zoltan_Migrate	Zoltan_Help_Migrate
ZOLTAN_PRE_MIGRATE_PP_FN	ZOLTAN_PRE_MIGRATE_FN
ZOLTAN_MID_MIGRATE_PP_FN	ZOLTAN_MID_MIGRATE_FN
ZOLTAN_POST_MIGRATE_PP_FN	ZOLTAN_POST_MIGRATE_FN

To continue using the v1.3 partition functions, no changes to C or Fortran90 applications are needed. Zoltan interfaces from versions earlier than 1.3 are also still supported (see [below](#)), requiring no changes to application programs.

To use the new v1.5 partitioning functions:

- C users must include file *zoltan.h* in their applications and edit their applications to use the appropriate new functions.
- Fortran90 users must put [user-defined data types](#) in *zoltan_user_data.f90* and edit their applications to use the appropriate new functions. The new partitioning functions do not work with user-defined data types in *lb_user_const.f90*.

Backward Compatibility: Versions 1.3 and higher

Versions of Zoltan before version 1.3 used a different naming convention for the Zoltan interface and query functions.

All functions in Zoltan v.1.3 and above are prefixed with **Zoltan_**; earlier versions were prefixed with **LB_**.

Zoltan versions 1.3 and above maintain backward compatibility with the earlier Zoltan interface. Thus, applications that used earlier versions of Zoltan can continue using Zoltan **without changing their source code**.

Only two changes are needed to build the application with Zoltan v.1.3 and higher:

- All Zoltan include files are now in directory *Zoltan/include*. Thus, application include paths must point to this directory.
(Previously, include files were in *Zoltan/lb*.)
- Applications link with Zoltan now by specifying only *-lzoltan*.
(Previously, applications had to link with *-lzoltan -lzoltan_comm -lzoltan_mem*.)

While it is not necessary for application developers to modify their source code to use Zoltan v.1.3 and above, those who want to update their source code should do the following in their application source files:

- Replace Zoltan calls and constants (**LB_***) with new names. The new names can be found through the [index below](#).
- C programs: Include file *zoltan.h* instead of *lbi_const.h*.
- F90 programs: Put [user-defined data types](#) in file *zoltan_user_data.f90* instead of *lb_user_const.f90*.

Backward Compatibility Index for Interface and Query Functions

Name in Earlier Zoltan Versions	Name in Zoltan Version 1.3 and higher
LB_BORDER_OBJ_LIST_FN	ZOLTAN_BORDER_OBJ_LIST_FN
LB_Balance	Zoltan_LB_Balance
LB_Box_Assign	Zoltan_LB_Box_Assign
LB_CHILD_LIST_FN	ZOLTAN_CHILD_LIST_FN
LB_CHILD_WEIGHT_FN	ZOLTAN_CHILD_WEIGHT_FN
LB_COARSE_OBJ_LIST_FN	ZOLTAN_COARSE_OBJ_LIST_FN
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