

MATCHER Manual

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1 Introduction

The prediction of many fluid/structure interaction phenomena requires solving simultaneously the coupled fluid and structural equations of equilibrium with an appropriate set of transmission (interface boundary) conditions. In many situations, the fluid and structure subproblems have different mesh resolution requirements and as a result, their computational domains have non-matching discrete interfaces ([Figure 1](#)). Converting the fluid pressure and stress fields at the fluid/structure interface into a structural load and transferring the structural motion to the fluid system can be performed in three steps:

- Step 1: properly pairing each “wet” structural element with a fluid entity (cell, control volume, or element), or an entity of the embedded discrete representation of the wet surface of the structure, as shown in [Figure 2](#).
- Step 2: computing the flow-induced structural loads, and structure-induced motion of the CFD grid points lying on the fluid/structure interface or the embedded discrete representation of the wet surface of the structure.
- Step 3: exchanging the resulting aerodynamic/hydrodynamic (loads) and elastodynamic (displacements and velocities) data between a flow solver and a structural analyzer.

In many applications, Step 1 can be performed once, in a preprocessing phase, by a computational geometry module. The **MATCHER** code discussed in this document is such a module. It outputs information that can be exploited to populate the data structures of a flow solver and a structural analyzer that are necessary for exchanging aerodynamic/hydrodynamic and elastodynamic data. Subsequently, Step 2 can be performed by elements of an accompanying library of subroutines that are callable from the flow solver and the structural analyzer. Finally, Step 3 can be performed either via any communication channel (for example, MPI) including I/O.

This document focuses on Step 1 and the **MATCHER** module.

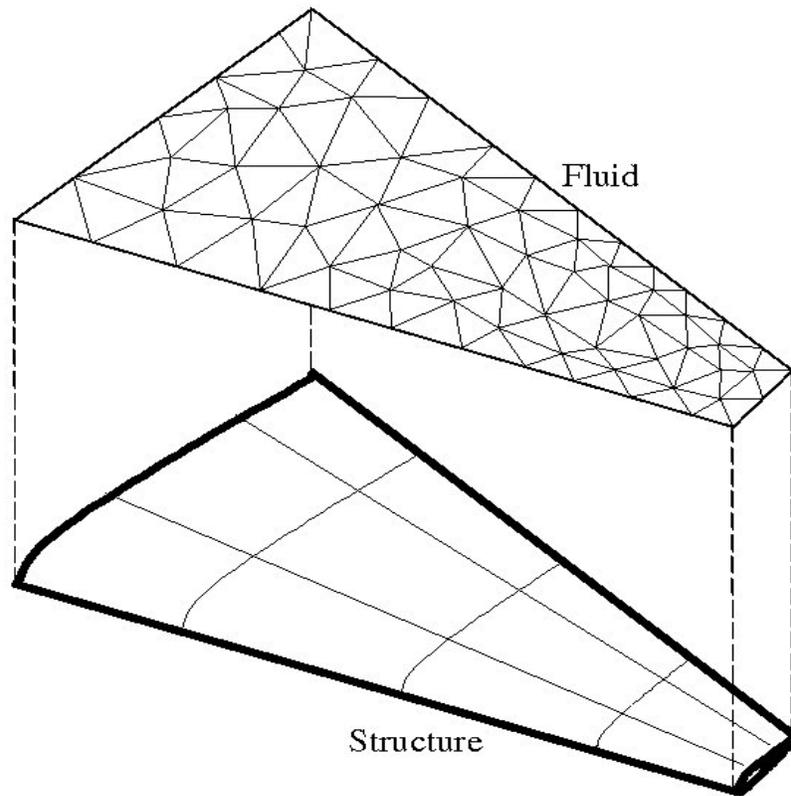


Figure 1: Fluid Mesh (or Embedded Discrete Representation of the Wet Surface of the Structure) and Structure Mesh with Non-Matching Discrete Interfaces

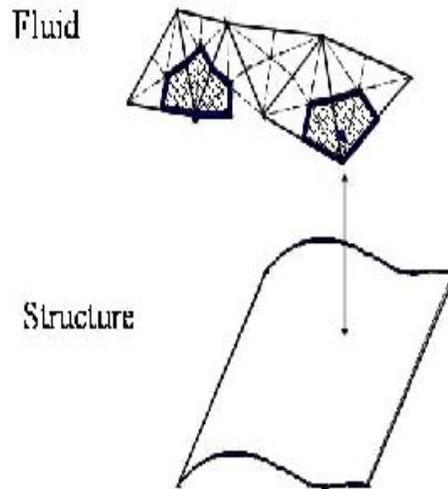


Figure 2: Pairing of Entities of the Structural Mesh and Fluid Mesh (or Embedded Discrete Representation of the Wet Surface of the Structure)

Next: [OutputFiles](#), Previous: [Introduction](#)

2 Input Files

MATCHER requires (and accepts) the following (optional) input files:

- `fluid.mesh`. This file, which can have any name, describes the unstructured mesh of the fluid domain. It can be in the **Sinus**, **XPost**, or **FieldView** format. When written in the **XPost** format, this file describes the unstructured discretization of the fluid domain as well as the boundary conditions. The latter are specified as element sets with the following conventions for naming the element sets: “OutletMoving” for the moving outflow boundary, “InletMoving” for moving inflow boundary, “StickMoving” for the moving viscous wall boundary, “SlipMoving” for the moving Euler wall boundary, “Symmetry” for the symmetry boundary, “SlipFixed” for the fixed Euler wall boundary, “StickFixed” for the fixed viscous wall boundary, “InletFixed” for the fixed inflow boundary, and “OutletFixed” for the fixed outflow boundary. Among these element sets, only the following ones are acted upon by **MATCHER**: “OutletMoving”, “InletMoving”,

“StickMoving”, and “SlipMoving”.

- `fluid.control`. This file, which can have any name, is an optional file. By default, only the nodes located on the “moving” surfaces of the fluid mesh are “matched” with corresponding nodes on the “wet” surfaces of the structural mesh. However, the user can request **either** that some nodes located on the moving surfaces of the fluid mesh be not matched, **or** that some nodes of the fluid mesh which are not located on the moving surfaces be matched. The unified syntax for both cases is given below.

```
CNBC < +- >                               <number of nodes>
<node number>
<node number>
...

```

A positive number of nodes on the first line indicates that the following nodes are to be matched with structure nodes, in addition to the wall surface nodes. A negative number of nodes on the first line indicates that the nodes specified on the following lines are not to be matched.

- `structure.input`. This file, which can have any name, is typically the **AERO-S** input file. Its content and syntax are described in the **AERO-S** User's Manual.
- `structure.control`. This file, which can have any name, is an optional file. By default, all the elements listed in the `structure.input` file are processed by the matching program. However, the user can request **either** that some elements listed in `structure.input` file be ignored, **or** that only a list of elements specified in the `structure.control` file be processed by the matching code. Furthermore, two different mechanisms based on elements and nodes, and which can be combined, are available for invoking **one or the other** action. The syntax for all cases is given below.

```
EIFS < +- >                               <number of elements>
<element number>
<element number>
...

```

```
NIFS < +- >                               <number of nodes>
<node number>
<node number>

```

...

A positive number of elements on the first line indicates that only the following structural elements are to be matched with fluid nodes. A negative number of elements indicates that the following structural elements are not to be processed even if listed in the `structure.input` file.

A positive number of nodes on the first line indicates that only the structural elements whose all nodes are among the following structural nodes are to be matched with fluid nodes. A negative number of nodes indicates that any structural element in the `structure.input` file which has any node among the following structural nodes is not to be processed by the matching code.

- `fluid.match`. This file, which can have any name, is an optional file. It can be used to prescribe the matching fluid nodes to structural elements on a one-on-one basis. The syntax for this file is given below.

```
<fluid_node_to_be_matched number>      <structural_element number>
<fluid_node_to_be_matched number>      <structural_element number>
```

...

- `structure.match`. This file, which can have any name, is an optional file. It can be used to prescribe the matching of structural elements to fluid nodes on a one-on-one basis. The syntax for this file is given below.

```
<structural_element_to_be_matched number>  <fluid_node number>
<structural_element_to_be_matched number>  <fluid_node number>
```

...

Next: [LineCommandInput](#), Previous: [InputFiles](#)

3 Output Files

The following output files are generated by **MATCHER** (**Warning: existing files are overwritten!**):

- `<prefix>.match.fem`. The detailed structure of this Structure MATCHER-NAME file is given in Appendix [StructureMatcher](#).
- `<prefix>.match.fluid`. The detailed structure of this Fluid MATCHER-NAME file is given in Appendix [StructureMatcher](#).

- `<prefix>.match.top`. This file contains the data needed for visualizing with **XPost** the outcome of the matching of the fluid and structure meshes.
- `<prefix>.nonMatch.top`. This file contains the data needed for visualizing with **XPost** all non matched fluid moving surface nodes.
- `<prefix>.bcdata`. This file, which is written in the **XPost** format of structural boundary conditions, contains all the non matched fluid moving surface nodes. The new convention adopted by **AERO-F** for non matched fluid moving surface nodes is that they are free to move. Hence, if the user wishes to restrain them, he/she can edit and/or paste this readily-available file into the fluid **XPost** input file to **MATCHER**.

Next: [SpecialFeatures](#), Previous: [OutputFiles](#)

4 Line Command Input

The line-command for running **MATCHER** in order to generate the auxiliary files needed by the **AERO-S** and **AERO-F** codes for performing a fluid/structure simulation is as follows:

```

matcher      <fluid.mesh file>
                <structure.input file>
                -b <fluid.control file>                (optional)
                -a <structure.control file>            (optional)
                -t <geometric tolerance>              (for computing intersection of normals and
                                                         elements; default value is 0.01)
                -e <distance tolerance>                (for canceling a performed matching when
                                                         the distance between the matched entities
                                                         exceeds the specified limit; default value is
                                                         MAXDOUBLE (a very large number))
                -m <fluid.match file>                  (optional)
                -n <structure.match file>              (optional)
                -beam <beam elements>                  (for processing beam elements)
                -l <number of levels for the geometric
                tolerance>                              (for repeatedly attempting to match
                                                         previously unmatched fluid elements)
                -output <output prefix>

```

-p <number of threads> (for running **MATCHER** in parallel mode using p threads)

In the above command, note that:

- The `fluid.mesh` and `structure.input` files must be input in the specified order. The optional features and their files can be specified in any order.
- The geometric tolerance is specified relative to the size of an element in its natural coordinate frame. Hence, it should be less than 1. The distance tolerance is specified in the absolute sense.
- When the option `-1` is specified, the geometric and distance tolerances are first set to their default or specified values. Then, repeatedly and as long as some fluid nodes remain unmatched and all specified levels of the tolerances have not been exhausted, the geometric and distance tolerances are multiplied by the level id number (1,2,3, ...) and the matching process is reattempted on the non matched nodes.
- The `-p` option can be used only when **MATCHER** is compiled with Intel/OpenMP.

Example:

- To generate the files needed for a fluid/structure analysis and excluding 11 fluid nodes from being matched

```
matcher          <fluid.mesh file>
                  <structure.input file>
                  -b <fluid.control file>
```

where the first line in the `fluid.control` file is

```
CNBC -11
```

- To generate the files needed for a fluid/structure analysis

```
matcher          <fluid.mesh file>
                  <structure.input file>
```

Next: [References](#), Previous: [LineCommandInput](#)

5 Special Features to Remember

- Beam elements in a 3D structural model are ignored by **MATCHER** unless the `-beam` flag is

invoked. Indeed, **MATCHER** assumes that when such beam elements are encountered on the wet surface of the structure, they are used to stiffen plate and shell elements that are also located on the wet surface of the structure and which share with these beam elements wet structural nodes. By default, **MATCHER** ignores these beam elements because their nodes can be found by processing the plate and shell elements that are also attached to them, and because the pressure induced finite element structural loads can be transmitted to these nodes by also processing only these plate and shell elements. However, if the `-beam` flag is invoked, **MATCHER** takes into account the beam elements of a 3D structural model and matches them with fluid nodes. (**This flag and therefore this option should be used only when there are no significant differentials in the motion between multiple beams — if the 3D structural model contains multiple beams — since neighboring points in the fluid mesh can take their input from neighboring beams**).

- **MATCHER** is parallelized for shared memory machines supporting the OpenMP standard. The number of threads is set via the environment variable, `OMP_NUM_THREADS`. For example, "setenv `OMP_NUM_THREADS 16`" will cause **MATCHER** to perform the interface matching with 16 threads, if possible. The default value is the minimum of 8 and the number of CPUs on the system.

Next: [ListofStructuralElements](#), Previous: [SpecialFeatures](#)

6 References

N. Maman and C. Farhat, *Matching Fluid and Structure Meshes for Aeroelastic Computations: A Parallel Approach*, Report CU-CUSSC-93-12.

C. Farhat, M. Lesoinne and P. LeTallec, *Load and Motion Transfer Algorithms for Fluid/Structure Interaction Problems with Non-Matching Discrete Interfaces: Momentum and Energy Conservation, Optimal Discretization and Application to Aeroelasticity*, Computer Methods in Applied Mechanics and Engineering, Vol. 157, pp. 95-114 (1998).

Next: [StructureMatcher](#), Previous: [References](#)

Appendix A List of Structural Finite Elements Supported by MATCHER

Element types	AERO-S Element type
• Bernoulli beam element	(6)
• Timoshenko beam element	(7)
• 3-node AQR shell element	(8)
• 4-node shell element element	(88)
• 3-node composite or orthotropic shell element	(20)
• 4-node composite or orthotropic shell element	(2020)
• 3-node triangular heat element	(46)
• 4-node triangular heat element	(4646)

Next: [FluidMatcher](#), Previous: [ListofStructuralElements](#)

Appendix B The Structure MATCHER File

This file is generated for fluid/structure simulations. It contains the data needed for relating the wet entities of the structure to the wall entities of the fluid. Currently, the structure is assumed to consist of a single global domain. An example of a **MATCHER** file is provided below.

```

SUBD 1 // number of structure subdomains
(always 1)
RCVF 0 // always 0
SNDF 2 // number of fluid subdomains w/
matched nodes
1 2 // fluid subdomain id & number of
matched nodes
741 0.23.069 0.277931 0.000000 0.000000 0.368834 // element number containing a matching
point, // natural coordinates of matching
point, // and gap vector
...
... // end of list
2 1 // fluid subdomain id & number of
matched nodes
719 0.197808 0.000000 0.000000 0.000000 -0.19567 // element number containing a matching
point, // natural coordinates of matching
point, // and gap vector
END SUBD // end of data for structural subdomain
END

```

Next: [GapVectors](#), Previous: [StructureMatcher](#)

Appendix C The Fluid MATCHER File

This file is generated for fluid/structure simulations. It contains the data needed for relating the wet entities of the structure to the wall entities of the fluid. An example of a this file is provided below.

```
1           // dummy subdomain id number (always 1)
1 3         // dummy number, number of matched nodes
2           // list of matched nodes
3
4
... ..     // end of list
0 0 -0.252963 // gap vectors
0 0 -0.268125
... ..     // end of list
```

Previous: [FluidMatcher](#)

Appendix D Gap Vectors

For non-matching fluid/structure meshes with non-coinciding geometric interfaces, there always exists gaps at the interface due to the difference in resolution of the fluid and structure meshes. This is shown in [Figure 3](#). These gaps must be considered when employing a mesh motion scheme in order to maintain the integrity of the meshes (i.e. prevent element interpenetration). For example, [Figure 4](#), shows the rotation of a flat plate when the gaps at the structure-fluid interface are not taken into account. The fluid nodes at the interface undergo the same displacements as the nodes at the structure interface. Thus, the application of a large rotation will eventually lead to element cross over. It is necessary to apply the rotation to the gap vector as well in order to maintain the mesh integrity as shown in [Figure 5](#).

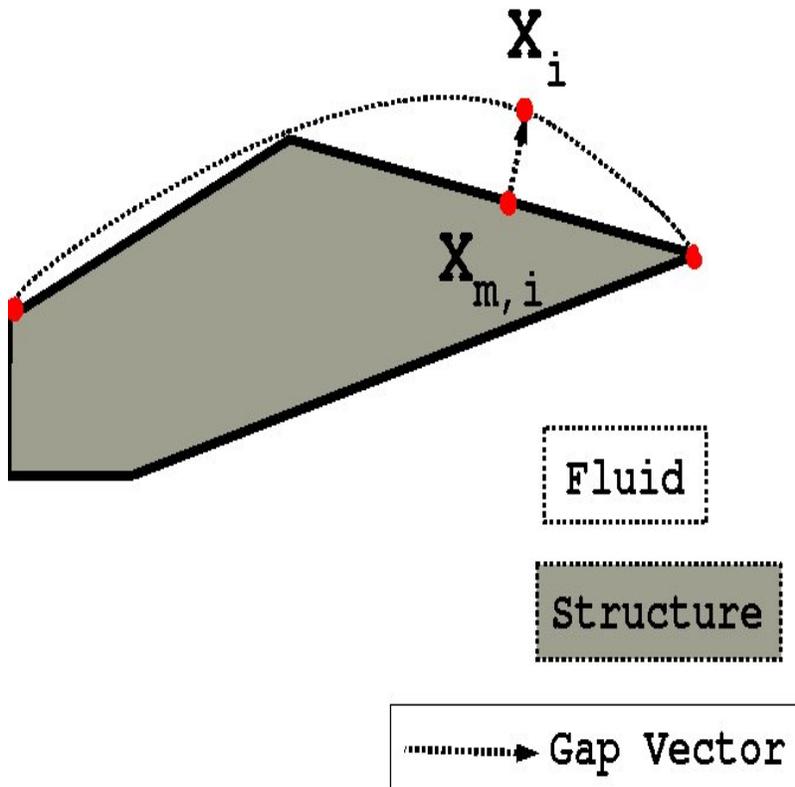


Figure 3: Fluid-Structure Interface

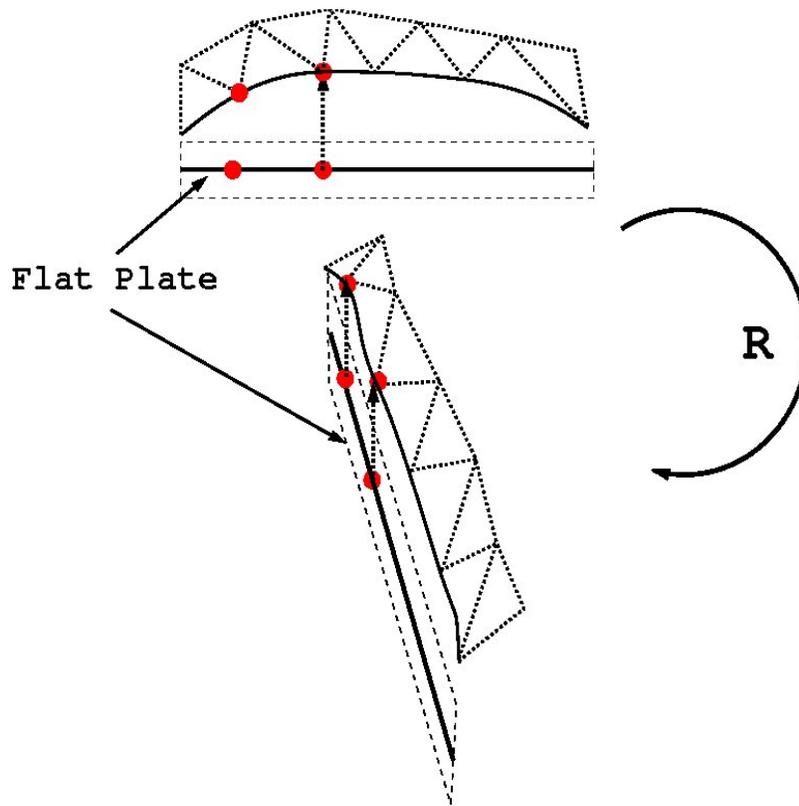


Figure 4: Flat Plate Rotation w/o Gap Vector

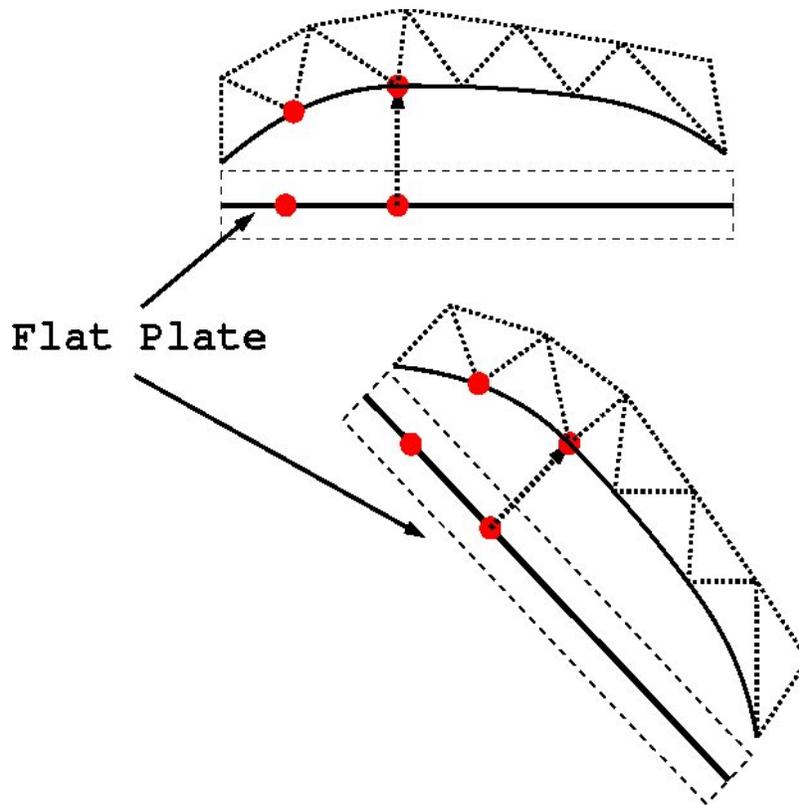


Figure 5: Flat Plate Rotation w/ Gap Vector