# **CFDEMcoupling Installation Guide**

Release 6.6.0

**DCS Computing GmbH** 

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## **A Product of**



 $CFDEM @ coupling \ stands \ for \ Computational \ Fluid \ Dynamics \ (CFD)-Discrete \ Element \ Method \ (DEM) \ coupling.$ 

CFDEM®coupling is an open-source code, distributed freely under the terms of the GNU Public License (GPL).

Core development of CFDEM®coupling is done by Christoph Goniva and Christoph Kloss, both at DCS Computing GmbH, 2012 (www.dcs-computing.com).

This documentation was written by DCS Computing GmbH, 2023.

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**CHAPTER** 

ONE

### INSTALLATION

This document will walk you through the installation process of CFDEM®coupling. First Aspherix® needs to be installed using its installer. OpenFOAM® and CFDEM®coupling are then installed. Finally, basic test cases are run to verify the successful installation.

As prerequisites OpenFOAM-10 and cmake >= 3.10 need to be installed. On Ubuntu 18.04+ cmake is installed during the OpenFOAM® installation with an appropriate version.

- Install Aspherix®
- Install and compile OpenFOAM®
- Install OpenFOAM® from binary package
- Install CFDEM®coupling
- Install CFDEM®coupling from binary package
- Test the installation
- Visualization using Paraview and Macro
- CFDEM Paraview reader
- Advanced installation information (optional)

## 1.1 Install Aspherix®

Before installing CFDEM®coupling, it is necessary that Aspherix® is propertly installed on your system. For more information regarding Aspherix®, please visit the official website. You receive details of the Aspherix® installation together with the Aspherix® installer as part of the provided documentation. The default installation location of Aspherix® on your system is \$HOME/DCS-Computing. You can test your Aspherix® installation by typing aspherix in your console. If successful, this command should create the output

Aspherix (Version ...) Checkout of asx\_solver OK.

### 1.2 Install and compile OpenFOAM®



The current version of CFDEM®coupling compiles *only* with OpenFOAM® in version 10.

If do not want to compile OpenFOAM® yourself but prefer to install from a pre- compiled package, refer to *Install OpenFOAM® from package*.

OpenFOAM is continuously updated and the source code repository contains numerous bugfixes that are not included in the pre-compiled package. To install and build the version that works best with CFDEM®coupling:

As a prerequisite for the OpenFOAM® installation you need to download and install additional software packages
(see the OpenFOAM webpage for more details). On Ubuntu 18.04 and newer execute the following command in
a terminal:

• Retrieve the source code. The following commands will create the required folders in an OpenFOAM directory in your HOME folder and download the desired version:

```
cd $HOME
mkdir OpenFOAM
cd OpenFOAM
git clone https://github.com/OpenFOAM/OpenFOAM-10.git
git clone https://github.com/OpenFOAM/ThirdParty-10.git
cd OpenFOAM-10
git checkout 362b7b6bb8eb5c4ac92e12856b3455cdacc1166a
cd ../ThirdParty-10
git checkout ae545a01baf66780285a2015da8a23c0a2dda5de
```

**Alternatively**, you can download the source code also as *zip archives*.

• "Enable" the version to be compiled and used by loading the OpenFOAM environment variables in your .bashrc file. Open your .bashrc file from the terminal (here using the default text editor in Ubuntu):

```
gedit ~/.bashrc
```

and add the following lines to the end of the file.

```
#- OpenFOAM SETUP
export WM_NCOMPPROCS=8
source $HOME/OpenFOAM/OpenFOAM-10/etc/bashrc
```

Here, 8 is the number of the processors you want to compile with, adjust this number to as many cores as your system has or you are willing to use. Save the changes.

• To compile, open a new terminal (!) so that above changes to your .bashrc become active and execute the following steps:

```
cd $WM_PROJECT_DIR
./Allwmake
```

Depending on your hardware and the amount of cores used, compilation of OpenFOAM can take an hour or more.

• For testing your installation, test that the simpleFoam application is found. Type into your terminal

simpleFoam -help

and execute it. You should get a "Usage" message printed to screen.

### 1.3 Install OpenFOAM® from binary package

Alternatively, you can install OpenFOAM® from the binary package provided by unpacking the archive to a suitable location.



### 1 Note

Depending on your specific Linux version, the binary package provided by the system or some external source may not provide a suitable OpenFOAM version. If you want to install a binary package, please use the package provided by DCS Computing.

### 1.4 Install CFDEM®coupling

You should have received a zip archive named "CFDEMcoupling-Treibacher-addTuts-6.6.0" containing CFDEM®coupling. Perform the following to install CFDEMcoupling:

• Unzip the archive into the folder \$HOME/DCS-Computing.



Do not unzip the zip archive in Windows directly, but do this under Linux or from within the WSL. Since the Windows file system is not case sensitive, unzipping the archive directly in Windows may lead to missing files and, consequently, errors during compilation.

• Set required environment variables and load the CFDEMcoupling bashrc. To do so, add the following lines your ~/.bashrc and make sure to add them *after* loading the OpenFOAM bashrc:

Setting the correct environment variables is crucial. In some cases you might need to deviate from the above template:

- If you are using a different Aspherix® version, e.g. a bugfix release, you may need to change the CFDEM\_ASX\_INSTALL\_DIR variable accordingly.
- If you use c-shell, manipulate ~/.cshrc accordingly.

### 1 Note

As a shortcut or as a blueprint you can use the sript load\_CFDEMcoupling-Treibacher-addTuts-6.6. 0.sh located at the root of the zip archive.

If you directly load the environment using this script and do *not* update your .bashrc, you need stay in the same terminal and load this environment in every new terminal you want to use CFDEM®coupling in.

### **A** Warning

The CFDEM bashrc sets many environment variables in a persistent manner. Those are **not** changed when sourcing that file a second time. Call cfdemUnsetEnv before changing the CFDEM environment or (re)-loading the CFDEM bashrc to ensure all variables are correctly updated.

- Open a new terminal so that above changes to your .bashrc become active
- Test your setup using the system test command:

#### cfdemSvsTest



#### 1 Note

Please note the output of the number of threads per core and the recommendation for the CFD-DEM hyperthreading setting.

· Compile using

### cfdemCompCFDEM

This command automatically compiles all the CFDEM®coupling libraries, CFDEM®coupling solvers and CF-DEM®coupling utilities. The compilation will stop, if build errors occur.

The compilation is automatically logged and the logs can be found in:

\$CFDEM\_SRC\_DIR/lagrangian/cfdemParticle/etc/log



#### 1 Note

If compiling on a different Linux distribution than Ubuntu, further settings may be required.

In case questions concerning the installation arise, please feel free to contact DCS Computing.



The above installation instructions, assume you have already installed Aspherix® in the default location \$HOME/DCS-Computing. Furthermore, we assume you have installed OpenFOAM®. If you are not using the CFDEM®coupling version CFDEMcoupling-Basic, please change the commands according to your CF-DEM®coupling version.

### 1.5 Install CFDEM®coupling from binary package

- Unpack the provided binary .tar.gz package to a suitable location. If you install from the .deb package, the default install location is /opt/CFDEM/CFDEMcoupling-Treibacher-addTuts-6.6.0.sh.
- Source the provided setup file to load the environment:

source ~/CFDEM/CFDEMcoupling-Treibacher-addTuts-6.6.0/load\_CFDEMcoupling-Treibacher-addTuts-6.  $\rightarrow$ 6.0.sh

If you unpacked to a different location, adjust the command accordingly. Preferably, add this command to your .bashrc.



#### 1 Note

The sourced script looks for OpenFOAM in /opt/OpenFOAM/ and \$HOME/OpenFOAM/. If you installed OpenFOAM in a different location, also update the OpenFOAM install path.

### 1.6 Test the installation

A simple way to test your installation is by running a tutorial case. In order to run a single tutorial it is recommended to use the Autorun.sh scripts in the tutorial directories (or their older counterparts Allrun.sh). They execute the simulation and provide some basic output (e.g., comparison of simulation output with an analytic solution).

Now, we describe briefly how to run a simple test case - cfdemSolverPiso/settling - which models the settling of a single sphere using cfdemSolverPiso and plots its velocity.

Open a terminal and navigate to the tutorial folder:

cfdemTut

cd cfdemSolverPiso/settlingTest



### 1 Note

If you do not load the environment in your .bashrc, you need to load it before the cfdem commands are available. Use, for instance, the sript load\_CFDEMcoupling-Treibacher-addTuts-6.6.0.sh located at the root of your CFDEMcoupling installation.

Run the simulation by executing the Autorun. sh script in the terminal:

./Autorun.sh

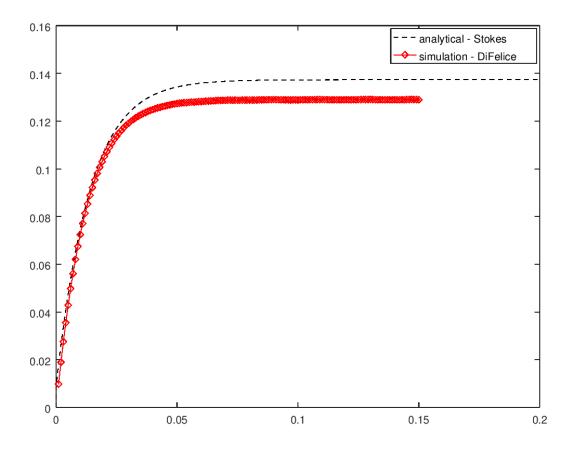
The executed script will prompt to select a variant for the test cases to run:

Available case variants (base variant is usually 'cfdemrun'):

- 1.) cfdemrun
- 2.) cfdemrun-serial
- 3.) cfdemrun-DEMbasedDrag

Which do you want to run? (Enter index):

Enter 1 to select the base variant cfdemrun. The other variants will change the case to use serial execution (2) or use a different drag law (for 3 the DEMbasedDrag). This will trigger Aspherix® to start the DEM simulation and CFDEM®coupling to start a coupled CFD-DEM simulation. Both simulations will communicate via socket and exchange the necessary data, e.g. particle position and velocity is communicated from DEM to CFD and the drag force is communicated from DEM to CFD. When the simulation is finished, a plot of the particle settling velocity pops up. The simulation result is compared to the analytical solution. You can visualize the simulation data using, e.g. Paraview.



In above plot the particle velocity (settling velocity) is ploted over time in red. The black line gives the analytical solution for the settling velocity using the Stokes drag law. Therefore, a slight offset between red and black line is to be expected.

You can run all the tutorial cases sequentially by executing the alias cfdemTestTUT. For running pure Aspherix® cases, you can use the aliases

```
cfdemAspherix inputScriptName cfdemAspherixPar inputScriptName nOfProcs
```

In the tutorials section you find more details on setting up and running a tutorials.

### 1 Note

If you want to run your own cases, it is recommended to first copy the tutorial case to \$CFDEM\_PROJECT\_USER\_DIR/run which is automatically being generated when you use the cfdemRun alias to go there for the first time. E.g., copy one of the tutorial cases there, and adapt it to your needs before running. This helps you keeping the tutorial in its original clean and working state.

### 1 Note

Refer to the CFDEMcoupling HTML documentation for all full documentation of models and solvers. The docu-

mentation is part of the Aspherix documentation.

### 1.7 Visualization using Paraview and Macro

To facilitate data loading a paraview macro is available to split loaded OpenFOAM / CFDEMcoupling data into separate blocks for internal field, patches, and –if present – particles.



The macro is automatically installed during compilation. This means that you need to perform the steps described below only if you want to use the Macro on a system (or for a user) where no CFDEMcoupling is compiled.

### 1.7.1 Installation

To use this macro copy it from its location to the Paraview Macro folder:

cp -r \$CFDEM\_ETC\_DIR/paraview/CFDEMcoupling\* <pv\_macro\_dir>/

where the pv\_macro\_dir is:

- on Linux:
- \$HOME/.config/ParaView/Macros'
- on Windows
  - %userprofile%\AppData\Roaming\ParaView\Macros

On Linux (assuming the above location of the Paraview config folder) you can also use:

cfdemInstallParaviewMacro

### 1.7.2 **Usage**

Load the CFD results by opening any \*.foam file in the CFD or CFD/system directories. After loading CFDEMcoupling simulation results you can run the CFDEMcoupling macro to split up the multiblock data, see Figure 2. Select all blocks when loading the data, i.e. also the boundaries and particles (if present), see Figure 1.

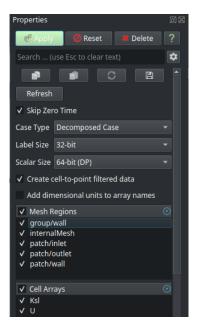


Fig. 1: Figure 1: Block selection.

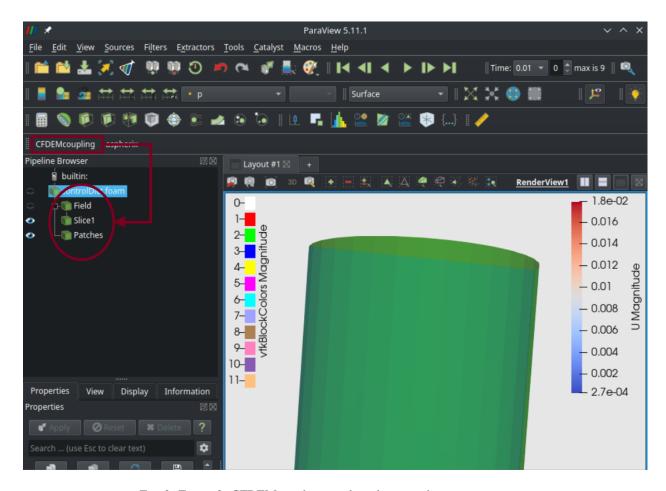


Fig. 2: Figure 2: CFDEMcoupling results split using the paraview macro.

### 1.7.3 Particle data

By default, particle data is written by Aspherix and stored in the DEM/post folder. To load all particle data, simply load the file aspherix\_simulation.pvd found in that directory.

To output particle data also on the CFD side, use an additional IOModel, e.g. basicIO.

### 1.8 CFDEM Paraview reader

In addition to the macro, a reader for coupled simulation results is installed automatically and should be available by default. If you install the macro manually as decribed above, the reader should also become available.

The CFDEMcoupling reader reads \*.cfdem files and exposes a node for, each, the CFD and the DEM results to which the corresponding macros may then attach. However, you still need to run the macros manually. The .cfdem file should be located in the root of the case directory.

### 1 Note

The reader is still at an early stage of development. If you encounter issues, please contact us.

### 1 Note

The usage of this reader is optional. You can always load the CFD and DEM results separately using the OpenFOAM and VTK-PVD readers.

## 1.9 Advanced installation information (optional)

In case the above instructions do not suit your needs, you can find advanced installation instructions in the following page:

### 1.9.1 Download OpenFOAM® zip archives

- Download the recommended OpenFOAM® version
- and the corresponding ThirdParty software.
- Unzip those archives to \$HOME/OpenFOAM and follow the other installation instructions.

### 1.9.2 Install OpenFOAM® from package



#### 1 Note

The current version of CFDEM®coupling compiles only with OpenFOAM® in version 10.

The easiest way to install OpenFOAM® on your system is by using the package manager to install a pre-compiled OpenFOAM package:

 As a prerequisite for the OpenFOAM® installation you need to download and install additional software packages (see the OpenFOAM webpage for more details). On Ubuntu 18.04 and newer execute the following command in a terminal:

```
sudo apt-get install build-essential cmake git ca-certificates flex libfl-dev bison.
→zlib1q-dev libboost-system-dev libboost-thread-dev libopenmpi-dev openmpi-bin
→ gnuplot libreadline-dev libncurses-dev libxt-dev
```

• Retrieve the OpenFOAM® package provided by OpenFOAM.org (find complete instructions for Ubuntu), which can easily be installed using the "apt" package manager. Basically the dl.openfoam.org repository must be added once to the list of software repositories. This is done with the terminal command:

```
sudo sh -c "wget -0 - http://dl.openfoam.org/gpg.key | apt-key add -"
sudo add-apt-repository http://dl.openfoam.org/ubuntu
```

• Install the downloaded OpenFOAM® package using the terminal commands:

```
sudo apt-get update
sudo apt-get -y install openfoam10
```

• "Enable" the installed package by loading the OpenFOAM environment variables in your .bashrc file. Open your .bashrc file from the terminal (here using the default text editor in Ubuntu):

```
gedit ~/.bashrc
```

and add the following line to the end of the file.

```
source /opt/openfoam10/etc/bashrc
```

Save the changes.

• For testing your installation, open a new terminal (!) and test that the simpleFoam application is found. Type into your terminal

```
simpleFoam -help
```

and execute it. You should get a "Usage" message printed to screen.

### 1.9.3 OpenFOAM® and MPI

#### 1 Note

This section discusses cases where a user-provided MPI is to be used with OpenFOAM. This is usually the case on clusters. On normal desktop systems, the default setting, i.e. export WM\_MPLIB=SYSTEMOPENMPI should work and does not need to be changed.

Often problems arise configuring OpenFOAM® to work with a specific MPI installation on a cluster. However it has built-in capability to work with nearly any MPI implementation. Set the following variable in the OpenFOAM® bashrc:

```
export WM_MPLIB=SYSTEMMPI
```

Defining the following variables in your ~/.bashrc before the sourcing of OpenFOAM® will take care of the rest. This is an example for older MVAPICH installations where there is no libmpi.so, only libmpich.so:

```
export MPI_ROOT=<path/to/mpi/installation>
export MPI_ARCH_PATH=$MPI_ROOT
export MPI_ARCH_FLAGS="-DMPICH_SKIP_MPICXX"
export MPI_ARCH_INC="-I$MPI_ARCH_PATH/include"
export MPI_ARCH_LIBS='-L$(MPI_ARCH_PATH)/lib -lmpich -lmpichcxx -lmpl -lopa -lrt'
```

### 1.9.4 Compiling CFDEM®coupling on different linux distributions

On some systems, like RHEL and CentOS, the MPICXX library is required in addition and you need to manually add it to the system linker paths *prior to compilation*. To do so, add

```
CFDEM_ADD_LIBS+=-lmpicxx
CFDEM_ADD_LIBS+=-L/path/to/lmpicxx
```

to

 $\verb|src/lagrangian/cfdemParticle/etc/addLibs_universal/additionalLibs_solver.|$ 

### 1.9.5 Compiling OpenFOAM® and CFDEM®coupling in debug mode

Set the  $WM_COMPILE_OPTION$  parameter in the OF bashrc to Debug and compile OF with the Allwmake script. CF-DEM®coupling detects the OpenFOAM® variable and compiles in debug mode.

### 1.9.6 Install swak4FOAM (optional)



Since migrating CFDEMcoupling to OpenFOAM-10, swak4FOAM is no longer used in the CFDEMcoupling tutorials. Consequently, the swak4FOAM installation is not tested with OpenFOAM versions beyond OpenFOAM-8.

swak4FOAM is a library of tools which open functionality otherwise only available by implementing new models. It is used in some test cases, mostly for post processing. To run these test cases without swak4FOAM, remove the related calls which are usually found in the functions subDict in controlDict.

To install swak4FOAM go to your OpenFOAM folder and run

```
# clone repo into swak4Foam sub folder
git clone https://github.com/Unofficial-Extend-Project-Mirror/openfoam-extend-swak4Foam-
dev.git swak4Foam

# go to cloned folder
cd swak4Foam
# checkout version compatible to OpenFOAM-8
git checkout 0f570f524c0e3be44155cdf0f7623c5403f234d3

# build
./AllwmakeAll
```

All dependencies and package requirements should already be installed from the package installation during the Open-FOAM installation process.



If encountering compilation problems with swak4FOAM, try first to run the compilation on a single core by running export  $WM_NCOMPPROCS=1$  before invoking ./AllwmakeAll.

### 1 Note

Refer to swak4Foam for further documentation on swak4FOAM.

### 1.9.7 Backward Compatibility

Basically, CFDEM®coupling supports one OpenFOAM® version. Therefore, all settings are prepared for that particular version. Nevertheless we try to maintain backwards compatibility as long as it works with reasonable effort.

The supported OpenFOAM® and Aspherix® versions are stated in:

 $\verb|src/lagrangian/cfdemParticle/cfdTools/versionInfo.H|\\$ 

For using other versions you can manipulate:

src/lagrangian/cfdemParticle/etc/OFversion/OFversion.H

### 1.9.8 Install git

Git allows you to update the source code very easily with the git pull command. On debian based systems you can run

```
sudo apt-get install git-core
```

Similar packages are available on other systems too:

```
sudo zypper install git-core
sudo yum install git
```

### **1** Note

In this tutorial the git protocol is used to transfer files. If port 9418 is closed for internet connections, please switch to "https://" instead of "git://" for git clone commands.

### 1.9.9 Install Helyx®OS (optional)

If you are interested in using a GUI based workflow for setting up your coupled CFD-DEM cases we recommend having a look into *Helyx*®-OS provided by Engys.



This is to be considered only a first "proof of concept" of combining CFDEM®coupling and Helyx®OS, which is planned to be streamlined in the future.

The detailed installation instructions for Helyx®OS basically boil down to downloading the installer from this link.

Then you need to change the permissions of the file by typing to the terminal

chmod 755 HELYX-OS-2.4.0-linux-x86\_64.bin

And finally you need to execute the installer in a terminal by typing

./HELYX-OS-2.4.0-linux-x86\_64.bin

### 1 Note

It is not required to install GUI + Kernel as you can use the OpenFOAM® version installed as described *above*. So please select only the components HELYX-OS-GUI. During installation the destination folder is being queried. As a default your user home directory is a good starting point. Simply type "echo \$HOME" in a terminal to see the path.

For a simple start of Helyx®OS we recommend to add to your ~/.bashrc (right above the definition of CF-DEM\_VERSION) the path of your Helyx®OS installation (you can find the correct path by looking for the file HELYX-OS.sh):

export CFDEM\_HELYXOS\_DIR=\$HOME/Engys/HELYX-OS/v2.4.0

In order to check your environment variables you can use the command

#### cfdemSysTest

, which will show your settings and if they are complete and correct. After defining CFDEM\_HELYXOS\_DIR you can start Helyx®OS by typing to your terminal:

### cfdemHelyxOS

After the first start you will have to define the paths to your OpenFOAM® installation. Please add

/opt/openfoam6

to Edit/Preferences/Core Folder.

### 1.9.10 Install OpenFOAM on Windows

You can run OpenFOAM® and CFDEM®coupling natively on Windows unsing the Windows sub-system for Linux (WSL).

1. To prepare your Windows system, check that you have updated your Windows 10 to at least version 2004, e.g. displayed in *Settings -> System -> Info -> Windows specifications*. If you do not have this version installed, run the update from *Settings -> Update & Security -> Windows Update*.

### 1 Note

You can force the update of Windows 10 to version 2004 as described in this manual.

2. Activate WSL by opening the PowerShell as Administrator via the start menu and execute the command:

```
{\tt Enable-WindowsOptionalFeature~-Online~-FeatureName~Microsoft-Windows-Subsystem-Linux}
```

When prompred choose to restart Windows.

3. When rebooted, open a new PowerShell terminal with Administrator privileges and execute:

```
Enable-WindowsOptionalFeature -Online -FeatureName VirtualMachinePlatform
```

Again, when prompted choose to restart Windows.

4. After the second reboot, start a new PowerShell terminal with Administrator privileges and to switch from WSL1 to the newer version WSL2, execute:

```
wsl --set-default-version 2
```

Now, WSL2 is set up and we proceed by installing a Ubuntu virtual machine which will be used by WSL.

- 5. Start by opening the Microsoft Store and search for Ubuntu 20.04. Click *Install*. This will download the system image and may take some time depending on your internet connection.
- 6. Start Ubuntu 20.04 from the start menu. This step will set up Ubuntu. In the process you must choose a username and a password for that user. These do not need to be identical with your Windows credentials.
- 7. Within the Ubuntu system run the following commands to update the package repository and install required tools:

```
sudo apt update
sudo apt install build-essential
```

8. OPTIONAL graphics support Install pre-requisites

```
sudo apt install nautilus gedit
sudo apt install xfce4
```

Open your bashrc (e.g. using nano)

```
nano .bashrc
```

and add the line

```
export DISPLAY=$(cat /etc/resolv.conf | grep nameserver | awk '{print $2}'):0
```

at the file end. Save the changes. If using nano, this is done using Ctrl + o followed by Ctrl + x.

9. Create links to acces the Windows file system more easily:

ln -s /mnt/c/Users/yourUserName/Downloads winDownloads

ln -s /mnt/c/Users/yourUserName/Documents winDocuments

- 10. Close the terminal.
- 11. Install Paraview on Windows
- 12. OPTIONAL graphics support

To run graphical programs on the Linux machine you need to additionally install an Xserver on your Windows. Fir this purpose we recommend VcXsrv. After installing VcXsrv, start it using the application Xlaunch.

1 Note

When doing so make sure to activate the checkbox "Disable access control".

1 Note

Modern systems may require to enable "Native opengl" to allow running the Aspherix-GUI.

13. OPTIONAL graphics support

Restart the ubuntu terminal.

In the terminal type

startxfce4 &

to start the graphical user interface.

Now, you are ready to proceed with the normal installation of *OpenFOAM* and *CFDEMcoupling* as described above within the just set up Ubuntu sub-system.

1 Note

After the installation of Aspherix-GUI you need to change the path of the Paraview installation to point to the one installed in Windows.

In case futher questions concerning the installation arise, please feel free to contact DCS Computing.							

### **1** Note

Refer to the CFDEMcoupling HTML documentation for all full documentation of models and solvers. The documentation is part of the Aspherix documentation.

#### **Footnotes:**

OPENFOAM® is a registered trade mark of OpenCFD Ltd, producer and distributor of the OpenFOAM software via https://www.openfoam.com.

HELYX® is a registered trade mark of Engys Limited, producer and distributor of the Helyx software via https://engys. com.