**User Manual**

|  |
| --- |
| **Compiling the source code** |

To compile the OPENDTS source code just execute “compile.sh” script.

./compile.sh

This will create three binary files: **DTS**, **CNV**, and **GEN**. The **DTS** is for running simulations; **GEN** is for generating triangulated files (TS) and **CNV** is for converting different file formats.

|  |
| --- |
| **How to run a simulation using OPENDTS** |

To run a simulation using **OPENDTS**, we need to feed at least two files to the **DTS** binary. An input file (with *dts* extension) that includes all the simulation inputs (see the “Input File” section) and a topology file that contains information about the topology of the TS mesh. The topology file could be in *top* or *tsi* format (see the Topology File sections).

With these two files a simulation can be started as.

$PATH/DTS -in input.dts -top top.top

Some of the parameters that are defined in the input file could be overwritten in the command line. Note: running a restart simulation is only possible by providing the restart file name in the command line as: -restart restart\_file\_name.res

|  |
| --- |
| **Simulation Output files** |

There are a number of output files for each DTS simulations that are listed below.

**vtu files:** This is a list of files that shows the evolution of the system and can be opened by paraview.

**tsi files:** A list of files that shows the evolution of the system. These files can be converted to different file format using the Convert Script.

**restart file:** A file to restart a simulation.This file has “res” extension.

**log file:** Stores some information of about the current simulation. This file has “*log”* extension.

**energy file:** This file end with *“-en.xvg*” and contains system the bending energy, box size, volume and etc., at each step.

|  |
| --- |
| **Command line options** |

All the option that can be defined in the command line. Also by executing .*/DTS -h* you can see this list.

|  |  |  |  |
| --- | --- | --- | --- |
| Identifier | Type | Default value | Description |
| -in | string | Input.dts | input file name |
| -top | string | topology.top | Topology file name |
| -b | int | 1 | initial time step |
| -e | int | 100 | final time step |
| -seed | int | 36723 | Random number seed |
| -defout | string | output | A general string for label a specific run |
| -ndx | string | Index.inx | Index file name |
| -restart | string | NO | restart file name |
| -angle | double | -0.5 | minimum of cos of the angle between two faces |
| -minDist | double | 1 | Square of minimum distance between two vertices |
| -maxDist | double | 3 | Square of maximum link length |

|  |
| --- |
| **Input file option and format** |

The input file must have a dts extension. Below, you can find the most common options defined in this file. However, this file could contain more information, including inclusions and their interactions. Other options are defined in their corresponding sections.

Integrator = MC

MC\_Moves = 1 1 1

Initial\_Step = 1

Final\_Step = 1000000

Display\_periodic = 1000

OutPutEnergy\_periodic = 1000

Restart\_periodic = 10000

Kappa = 15 0

Spont\_C = 0

Seed = 37382

OutPutTRJ\_TSI = 1000 10 TrjTSI

GeneralOutputFilename = output

Cell\_Size = 2.5 2.6 2.5

**Integrator:** The algorithm for updating the configuration of the system. Currently only “MC”available.

**MC\_Moves:** An option for activating or disactivating the MC moves.First one is vertex move, second is the link flip and the last one is the inclusion moves. 1 1 1 means all the moves are active.

**Initial\_Step:** Initial step of the simulation, usually it should be one, unless something else is intended.

**Final\_Step:** The final step of the simulations

**Display\_periodic**: Frequency of coordinate file in vtu file format.

**OutPutEnergy\_periodic:** Frequency of energy file. Energy file contains several other information depending on global constrains applied on the system.

**Restart\_periodic**: Frequency of restart file recording.

**Kappa:** Membrane bending and gaussian rigidity.

**Spont\_C:** spontaneous curvature of the membrane

**Seed:** Random number seed.

**OutPutTRJ\_TSI:** Frequency for writing trajectory frames in tsi file, the precision of the coordinates and the name of the folder that the files will be stored in.

**GeneralOutputFilename:** A general string to label all the generated files of a specific run with that string.

**Cell\_Size**: The size of the unit cells for domain decompositions, should be larger than 1

**Box\_Centering\_F:** To canter the system inside a box after every m steps. The m should be provided as an input. This

**FreezingAGroup:** Freezes a group of vertices.

FreezingAGroup

MinfaceAngle

TopologyFile

CouplingtoFixedGlobalCurvature

Min\_Max\_LinkLenghtsSquare

Parallel\_Tempering

OutPutTRJ\_BTS

Mem\_model\_para

**MC\_MovesRate this has been deactivated in the code**

**Fixed frame tension simulations:** For doing this, one needs to add below line to the input file. The last two numbers are the frame tension in the unit of [kBT/a2] and inverse frequency of applying the box size check algorithm to maintain the fixed membrane tension respectively. Note: this algorithm only makes sense for predoc membranes.

Frame\_Tension = on Position\_Rescale 0 5

**Pressure deference or fixed volume simulation.**

;Volume\_Constraint = on eqtime DP K gamma

Volume\_Constraint = on 100 -0.1 0 0

**Apply Constant Area**

**Applying harmonic potential on the surface global curvature.**

;CouplingtoFixedGlobalCurvature = state K gC0

CouplingtoFixedGlobalCurvature = on 60 0.1

**Membranes in confined spaces:** OPENDTS also allows simulating membranes while they are in confined regions. If the initial membrane is not within the defined region, bthe surface is forced to enter this region by defining an equilibrium time. However, if the eq time is short, it may fail to do so. There are four types of confided space defined in OPENDTS: TwoFlatParallelWall, Cuboid, Ellipsoid and EllipsoidalShell. To apply one of these confinements, one of the below commands must be added to the input file.

CoupleToRigidWalls = on TwoFlatParallelWall 10000 2

CoupleToRigidWalls = on Cuboid 10000 10 10 10

CoupleToRigidWalls = on Ellipsoid 10000 10 10 10

CoupleToRigidWalls = on EllipsoidalShell 10000 10 10 10 0.01

**Harmonic potential between two groups of vertices**: To apply a harmonic potential between two groups of vertices, one need to add below command to the input file.

; HarmonicPotentialBetweenTwoGroups = on K L0 eqstep Group1 Group2 nx ny nz

HarmonicPotentialBetweenTwoGroups = on 10 100 2000000 Group1 Group2 0 0 1

You also need to provide an index file for the group definition. This file should have “*inx”* extension. An example of index file is as below:

Group1 1

24

Group2 3

12 16 33

**Adding inclusions:** To have inclusion in the simulations, we need to define different type of inclusions. Inclusion type must be defined at the end of the input file as

INCLUSION

Define 4 Inclusions

SRotation Type K KG KP KL C0 C0P C0L

3 Pro1 10 0 0 0 0 0 0

3 Pro2 20 0 0 0 0 1 0

3 Pro3 20 5 0 0 0 0 0

2 Pro4 20 5 0 0 0 0 0

The Inclusion-inclusion interactions should be always placed at the end of the file as

Inclusion-Inclusion-Int

1 1 1 2 0.0 0.0

Type 1 n A B

Defining inclusions type inside the input file does not means that they are present in the simulation. To create inclusion in the simulations, we need to define them either inside the “*tsi*” file or tell the input file to generate them. Currently we can only generate random distribution of inclusions from the input file. However, some can manually or use a script to add such inclusions to the tsi file.

GenerateInclusions

Selection\_Type Random

TypeID 1 2 3

Density 0.0 0 0

|  |
| --- |
| **Osmotic pressure** |

Membranes are impermeable to ions and large molecules.

Jacobus van 't Hoff found a quantitative relationship between osmotic pressure and solute concentration

is the dimensionless van 't Hoff index and is the molar concentration of solute. For most non-electrolytes dissolved in water, the van 't Hoff factor is essentially 1.

For a vesicle with a volume, we have

In the code:

Osmotic\_Pressure = Type time gamma P0

Where

|  |
| --- |
| **Topology file format** |

Topology file provides information about the position of all the vertices and how they are linked. There are two types of file that you can provide as a topology file; *top* and *tsi* file format.

***top* file:** This file has a “top” extension and contains a list of TS files in q (TS file) file format. An advantage of this topology format is that it allows us to feed multiple TS file into the system. A disadvantage is that it does not include inclusions.

***tsi* file**: tsi topology is single file that is also single frames of OPENDTS trajectories. An advantage of this topology file is that it includes the inclusions information.

**TS file**

Triangulated surface files that can be read by OPENDTS either has “q” format or “tsi” format. The Generate script can generate files in both formats (see Generate script section).

**“tsi” format files:**

The following shows a part of a .tsi file with all necessary keywords highlighted in bold. Every .tsi file starts with a line calling version 1.1. The next line defines the box size (x, y, and z) of the system in nm. The next three sections describe the TS mesh. Each section starts with a keyword (vertex, triangle and inclusion) and their corresponding number. Here, we have 130 vertices (the numbering starts from 0). Each vertex has an index and a position in x, y and z (in nm). The 130 vertices are connected via 256 triangles. Again, every triangle has an index (starting from 0) and is defined by the vertices the triangle connects, i.e. triangle 0 connects vertices 11, 55 and 43. Furthermore, a .tsi file can have a (protein) inclusion section. Here, there are three inclusions from two different types. Again, each inclusion has an index. The index is followed by the inclusion type (here: type 1 for inclusions 0 and 1, type 2 for inclusion 2) and the corresponding vertex index. The last two (floating point) numbers describe a unit two-dimensional vector (sum of both numbers must be one!) which defines the orientation of the inclusion with respect to the bilayer normal.

**version 1.1**  
**box**   50.0000000000     50.0000000000     50.0000000000  
**vertex**130  
0     21.1606233083     25.4394806652     25.5960855271  
1     27.0284995400     23.2012757654     21.6715285158       
2     26.9921761232     25.5136587223     28.0195776981  
3     23.3273229896     26.2315165676     28.0075875808       
4     26.2722773116     26.3271061222     28.1420707299  
5     22.0396876425     23.6080597437     26.8858740866       
.  
.  
.  
125   21.5556280860     25.5595098219     26.5363425272  
126   23.2182025326     26.8060871266     21.5195141902       
127   25.3199303865     24.3519379911     20.6752314764       
128   28.0093200458     22.6356946990     23.4685318698  
129   21.4000741257     26.5841316766     25.2761757772  
**triangle**256  
0     11     55     43  
1     94     75     14  
2     64      3     91  
3     59     52     40  
.  
.  
.  
253   33    109     44  
254   53     69     47  
255   85      6     74  
**inclusion**        3  
0      1     22      0     1  
1      1      5      0     1  
2      2     30      0     1

**“q” format files:** The previous tsi file in q format can be seen below.

Line 1: Box information (3 double numbers).

Line 2: Number of vertices (1 integer number; Let’s call it NV).

Line 3 to NV+2: Vertex ID and coordinate (1 integer number and 3 double numbers).

Line NV+3: Number of triangles (1 integer number; Let’s call it NT).

Line NV+4 to NV+NT+3: Triangle ID and ID of its vertices (4 integer number).

50.0000000000     50.0000000000     50.0000000000  
130  
0     21.1606233083     25.4394806652     25.5960855271  
1     27.0284995400     23.2012757654     21.6715285158       
2     26.9921761232     25.5136587223     28.0195776981  
3     23.3273229896     26.2315165676     28.0075875808       
4     26.2722773116     26.3271061222     28.1420707299  
5     22.0396876425     23.6080597437     26.8858740866       
.  
.  
.  
125   21.5556280860     25.5595098219     26.5363425272  
126   23.2182025326     26.8060871266     21.5195141902       
127   25.3199303865     24.3519379911     20.6752314764       
128   28.0093200458     22.6356946990     23.4685318698  
129   21.4000741257     26.5841316766     25.2761757772  
256  
0     11     55     43  
1     94     75     14  
2     64      3     91  
3     59     52     40  
.  
.  
.  
253   33    109     44  
254   53     69     47  
255   85      6     74

|  |
| --- |
| **Generate Script** |

Generate (**GEN**) bindery allows you to create triangulated surface files with different topologies in two different file formats, *“q”* or *“tsi”*. Three options exist, flat bilayer periodic in x and y directions, cylinder periodic in the x-direction and a closed sphere. To see all the options, execute $path/GEN -h

Some example:

To generate flat bilayers

$path/GEN -box 50 50 30 -type flat -o topol.q

Generating closed membranes

$PATH/GEN -box 50 50 50 -type tetrahedron -N 20 -o topol.q

Note: You can force this structure to become sphere by using Ellipspoidalshell command and a short simulation.

CoupleToRigidWalls = on EllipsoidalShell 10000 10 10 10 0.01

|  |
| --- |
| **Convert Script** |

The convert (**CNV**) bindery allows for converting “*tsi”* and “*q*” files to each other and several other different file format such as “*vtu”* and “*gro*”. To see all the options, execute $path/GEN -h

**Tutorials**

***T1: Framed membranes***

The simplest simulation that can be performed by FreeDTS is a flat membrane (a fluid elastic surface) in a PBC box. To do this, we have to first generate a flat TS file. This can be done by using **GEN** binary. Using this you can generate a TS file in a q file format using the below command line.

$path/GEN -box 30 30 100 -type flat -o topol.q

This command creates a TS file in *q* file format with a box size of 50\*50\*100. Next, this file name should be added to a topology file name with top extension (here it is named top.top), and we should add the below line in the top file.

topol.q 10

The number in front of the file gives an *id* to the entire mesh (the value is not important but needs to be unique if multiple files are included). To run a simulation, an input file (with dts extension) is required to define the simulation parameters (see below box and the input file format section).

Integrator = MC

MC\_Moves = 1 1 1

Initial\_Step = 1

Final\_Step = 100000

Display\_periodic = 1000

OutPutEnergy\_periodic = 100

Restart\_periodic = 10000

Kappa = 20 0

OutPutTRJ\_TSI = 1000 10 TrjTSI

GeneralOutputFilename = output

Using these two files we are now able to run DTS simulations as:

$PATH/DTS -in input.dts -top top.top -seed 76532

Everything above could be done by running the *./run.sh* script in the dts\_tutorials/T1-FramedMembrane\_a folder in the source code.

To make a membrane tensionless, we need to add below command to the input file (such the example in the input file section).

Frame\_Tension = on Position\_Rescale 0 5

In the dts\_tutorials/T1-FramedMembrane\_b folder run the *./run.sh* script to perform this section.

Each run will give multiple outputs. Folder VTU\_Frames contains *paraview* readable files to visualize the evolution of the system. TrjTSI folder provides files in *tsi* format for analysis of the simulation. *output-en.xvg* file contains information about system energy, box size (for constant tensions simulations) membrane volume for constant volume simulations (see the vesicle tutorials) …. as a function of the simulation steps.

***T2: Vesicle simulations***

**Perpetrating a Vesicle**

For a simulation of a vesicle, we first need to create a vesicle structure. For this, we use **GEN** script to generate a tetrahedron and run a short simulation with *EllipsoidalShell* confinement to shape the tetrahedron into a vesicle. The below command creates a *TS* file in the shape of tetrahedron. The size of the tetrahedron can be changed by -N option

$PATH/GEN -box 50 50 50 -type tetrahedron -N 20 -o topol.q

Next, we add the generated TS file to the topology file and run a normal simulation while having below line in the input file (for more see “Membranes in confined spaces” section in the manual).

CoupleToRigidWalls = on EllipsoidalShell 10000 10 10 10 0.01

After about 20K steps, you should get a spherical TS file (the run steps may differ for different system)

In the dts\_tutorials/*T3-Vesicle* folder run the *./run.sh* script to perform this tutorial.

***Note****: one can easily create a TS file with a spherical shape. However, we prefer this method because it allow us to create a TS file with regular edge size (between range required for DTS simulations) on many different shape.*

**

**Constant Volume Vesicle**

Using the obtain spherical TS file, we can obtain a shape of a vesicle with a constant volume. For this use the **CNV** script and convert last frame of the previous simulation into a q file for a new simulation.

$PATH/CNV -in output20.tsi -o vesicle.q

Remove below line from your input file

CoupleToRigidWalls = on EllipsoidalShell 10000 10 10 10 0.01

And add below line to the input file (see “Osmotic pressure or fixed volume simulation” section)

Volume\_Constraint = on 10000 0 0.1 0.4

***T3: Proteins***

By adding below section to the end of the input file from the previous tutorial, we can add inclusions to the system (For more information, see the Input file section in the User Manual).

INCLUSION

Define 4 Inclusions

SRotation Type K KG KP KL C0 C0P C0L

0 Pro1 10 0 0 0 0.4 0 0

0 Pro2 20 0 0 0 -0.4 0 0

GenerateInclusions

Selection\_Type Random

TypeID 1 2 3

Density 0.3 0.1 0

Inclusion-Inclusion-Int

1 1 1 2 2 0.0

1 2 1 2 0 0.0

2 2 1 2 2 0.0

In the dts\_tutorials/*T2-Membranes\_Inclusions* folder run the *./run.sh* script to perform this tutorial.

***T4:* Pulling a membrane nanotube**

To pull a nanotube from a membrane, we need to have a membrane under tension (see previous tutorials for such a simulations). Then we can apply a harmonic potential between two groups, one group can be the single vertex while the other group can be all the other vertices. Using a corresponding index file (See the index file section) and adding below command to the input file we have everything we need to pull a nanotube from a flat membrane.

HarmonicPotentialBetweenTwoGroups = on 10 100 2000000 Group1 Group2 0 0 1

Next, just run below command and wait for the nanotube to form.

$PATH/DTS -in input.dts -top top.top -seed 76532 -ndx index.inx

In the dts\_tutorials/*T4-tether\_pulling* folder run the *./run.sh* script to perform this tutorial.

***T5:* Protein Sorting**

***T5:* two TS files system**