# VULTURE FDTD CODE

# User Manual

Vulture Version 0.7.0 Mesh Version 1.0.0 (9/12/2017)

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# Acknowledgements

• The input mesh format for Vulture is an evolution of the format used by the University of York "Hawk" Transmission Line Matrix and "Falcon" Finite-Difference Time-Domain solvers by Dr Stuart Porter and Dr John Dawson [\[1,](#page-65-0) [2\]](#page-65-1). The binary output format and the associated post-processing tools were also originally developed by Dr Stuart Porter and Dr John Dawson for these two codes [\[3\]](#page-66-0).

# **History**

# Contents





# <span id="page-4-0"></span>Acronyms and abbreviations



# <span id="page-5-0"></span>1 Introduction

# <span id="page-5-1"></span>1.1 Overview

This user guide provides both a user manual and tutorial introduction for the Vulture finite-difference time-domain (FDTD) code. Details about the theory and implementation of the code can be found in the Vulture Code Implementation Manual [\[4\]](#page-66-1). This user guide assumes that the Vulture software has already been installed according to the instructions in the binary or source distribution package as appropriate. Please also refer to the ReadMe.txt and Bugs.txt files in the distribution package for details of any known bugs or issues with the particular version of Vulture you have installed.

# <span id="page-5-2"></span>1.2 Solver features

Vulture is a non-uniform structured mesh FDTD code which has the following features:

- Non-uniform mesh allowing for uniform cubic and uniform cuboid special cases;
- External mesh surfaces that can independently be perfect electric conductor (PEC), perfect magnetic conductor (PMC), perfectly match layer (PML), analytic Mur absorbing boundary condition (ABC) or periodic boundary conditions;
- A uniaxial perfectly matched layer (UPML) implementation that can terminate arbitrary inhomogeneous media;
- Gaussian pulse, compact pulse, ramped sinusoid, differentiated pulse and user defined waveforms;
- Distributed hard and soft electric and magnetic field, current density, current and ideal voltage sources;
- Lumped resistive voltage and current sources;
- Internal PEC surfaces;
- Simple isotropic media with frequency independent permittivity, conductivity and (real) permeability.
- Arbitrary electrically dispersive media using a generalised multi-pole Debye dispersion relationship.
- A total-field-scattered-field (TFSF) plane-wave source, also known as a Huygen's surface source, for multiple plane-wave excitation. The implementation supports partial Huygen's surfaces and has grid dispersion optimisations for uniform cubic meshes.
- Binary and ASCII format field observers.

Additional optional features that can be enabled at compile time are:

- (WITH SIBC=ON/OFF): Enables/disables frequency dependent surface impedance boundary condition (SIBC) surface material types.
- (USE AVERAGED MEDIA=ON/OFF): Enables/disables volumetric material averaging. When enabled this providies second-order accurate treatment of boundaries between different media (incompatible with USE\_INDEXED\_MEDIA=ON, EXPERIMENTAL).
- (WITH OPENMP=ON/OFF): Enables/disables multi-threaded parallelisation of the core update algorithms for increased performance on shared memory multi-core computers.
- (USE INDEXED MEDIA=ON/OFF): Enables/disables the use of indexed media. Indexed media can considerably reduced memory consumption, with a small penalty in run-time performance, providing the number of media is not very large (incompatible with USE AVERAGED MEDIA=ON and USE SCALED FIELDS=ON.
- (USE SCALED FIELDS=ON/OFF): Enables/disables the use of scaled fields on the computational grid. Scaled field give a small improvement in performance by reducing the operataion count in the discete curl operator (incompatible with USE INDEXED MEDIA=ON).

# <span id="page-6-0"></span>1.3 Running the solver

Vulture is a command-line program that runs from the Linux/Unix shell or Windows command prompt. For example, under Linux the program can be run using

\$ vulture -h

Usage:

```
vulture -h | --help
vulture -V | --version
vulture [ option ] <meshFile>
```
Valid options are:



where the  $-h$  option is used to provide basic usage information. The settings of the compilation options used to compile the executable can be checked using the -V option:

### \$ vulture -V

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Vulture comes with ABSOLUTELY NO WARRANTY; for details type 'vulture --licence'. This is free software, and you are welcome to redistribute it under certain conditions; type 'vulture --licence' for details.

Supported mesh versions 1.0.0 - 1.0.0 Using un-scaled fields. Field arrays are 4-bytes. Using indexed media. Medium index is 4-bytes. Using un-averaged media.

The program expects an input mesh file in the current directory which is in the format given in Section [1.4.](#page-7-0) The program's outputs are written into files in the same directory, hence each simulation requires its own working directory in order to avoid confusion and overwriting of output data. A typical command line for running Vulture is

# \$ vulture -v antenna.mesh

where  $antenna$  is the name of the input mesh file and the  $-v$  option requests more verbose information in the log file. The output files created by the solver include:

- vulture.log: An ASCII log file containing information on the mesh and simulation.
- xlines.dat, ylines.dat and zlines.dat: ASCII files containing lists of the  $x, y$  and  $z$  mesh line coordinates.
- impulse.dat: A binary format file containing time-domain field outputs.
- excite.dat: An ASCII file containing the excitation waveform time series.
- process.dat: A template ASCII file for controlling other post-processing tools.
- wf\_<t: name>\_td.asc: ASCII data files containing waveform time series.
- eh\_<t: name>\_td.asc: ASCII data files containing field time series.
- wf <t: name> fd.asc: ASCII data files containing waveform frequency spectra.
- eh <t: name> fd.asc: ASCII data files containing field frequency spectra.

If the solver has been compiled with OpenMP parallelisation support the number of threads used can be set wtih the command line flag -n. For example, to use two threads start the solver with

### \$ vulture -n 2 antenna.mesh

If the number of threads is not specified using the option argument then the value of the environment variable OMP NUM THREADS is used. On Linux/Unix system using Bourne/bash shells this can be set with

#### \$ export OMP NUM THREADS=2

If neither the option argument or environment variable are set then a single thread is used.

# <span id="page-7-0"></span>1.4 The mesh file

The simulation is defined in an ASCII mesh file with the format shown in Table [1,](#page-8-0) in which each line performs a specific function. The first two characters of each line constitute a two letter code or directive that specifies the nature of the information on the rest of the line. The # character can be used to insert comments anywhere in the mesh file, either full lines or at the end of a line. In the table it has been used to insert comments identifying the different logical parts of the mesh file. Empty lines can also be placed anywhere in the file. Each directive must appear on a single line in the file. In this manual some directives with long lines have been wrapped onto a second line denoted by the back-slash at the end of the first line. In the real input file the directives must be place on a single line. Optional elements in the directives are enclosed in square brackets [].

The VM directive must be the first directive in the mesh file. It specifies the version of the mesh format and is used by Vulture to determine how to parse the file. The other directives can occur in any order within their respective mesh sections. The mesh sections must be in the order shown in Table [1.](#page-8-0) The number of times each directive can be used (its multiplicity) and other constraints are listed in Table [2.](#page-9-0) Some directives are optional (multiplicity 0,1), with default values being assigned in their absence, while others must occur once and only once (multiplicity 1). Sections 1 and 3 of the mesh file define the computational grid while section 2 defines the objects on the grid.

The elements following the two-letter directive code are user defined values of the following types:

- <i: var>: An integer.
- <r: var>: A real number.
- $\leq$  s: var>: A general character string up to 1023 characters long. The string must be enclosed in double quotation marks, for example "waveform.asc".
- $\leq t$ : var>: A "tag", which is defined as a character string consisting of the letters "a-z", "A-Z". numbers "0-9" and underscore " $\cdot$ " with length up to 31 characters. The string must not be enclosed in quotation marks and should not start with an underscore character. These tags are used to name and reference entities in the mesh and must be defined before they are used.

The detailed meanings of all the mesh directives are given in Sections [2](#page-10-0) to [5](#page-36-0) below.

### <span id="page-7-1"></span>1.5 The log file

Vulture creates a detailed log file called vulture.log in the current directory. This file contains:

- Information about the parsing of the mesh file.
- A report on the input mesh characteristics.
- Initialisation information for all the grid objects.
- A report on the computational grid characteristics.
- A breakdown of the memory usage.
- A periodically updated estimate of the completion time of the simulation.

```
VM <i: majorVersion>.<i: minorVersion>.<i: patchVersion>
#
# Section 1
#
CE <s: comment>
DM <i: nx> <i: ny> <i: nz>
GS
#
# Section 2
#
MT <t: name> <t: type> <param1> ... <paramN>
BT <t: name> <t: type> <param1> ... <paramN>
WT <t: name> <t: type> <param1> ... <paramN>
MB <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> [ <t: mask> ]
TB <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> [ <i: orient> [ <r: angle> ] ]
TW <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> [ <t: end1Type>] [ <t: end2Type> ]
WF <t: name> <t: type> [ <r: size> [ <r: delay> [ <r: width> [ <r:freq> ] ] ] ]
WF <t: name> EXTERNAL <s: fileName> [ <r: size> [ <r: delay> ] ]
EX <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: type> <t: wfName> \
   [ <r: size> [ <r: delay> ] ]
EX <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: type> <t: wfName> \
   <r: imped> [ <r: size> [ <r: delay> ] ]
PW <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: wfName> \
   <r: theta> <r: phi> <r: psi> [ <i:mask> [<r: size> [ <r: delay> ] ] ]
OP <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: type> [<t: wfName>]
OP <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: type> [ <t: wfName> \
   [ <i: xstep> <i: ystep> <i: zstep> ] ]
FF <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: wfName> \
   <r: theta1> <r: theta2> <i: num_theta> <r: phi1> <r: phi2> <i: num_phi> [ <t: mask> ]
GE
#
# Section 3
#
NT <i: numSteps>
CN <r: courantNumber>
OT <i: tstart> <i: tstop>
OF <r: fstart> <r: fstop> <i: numFreq>
MS <r: delx> [ <r: dely> <r: delz> ] ]
Y<sub>I</sub><r: x[0]>
.....
\langle r: x[nx] \rangleYL
<r: y[0]>
.....
<r: y[ny]>
71
\langle r: z[0] \rangle.....
\langle r: z[nz] \rangleEN
```
<span id="page-8-0"></span>Table 1: Input mesh file format. Note that some directives with long lines have been wrapped onto a second line denoted by the back-slash  $(\rangle)$  at the end of the first line. In the real input file the directives must be place on a single line. Optional elements are enclosed in square brackets [].

• A final summary of the run-time and the overall average seconds per iterations (spi) per cell.

If called with the -v option Vulture will add more detailed information to the log file that is useful for debugging the solver and identifying problems with the mesh.

An example of the grid report from the log file is:

```
Grid characteristics:
```

```
Grid is CUBIC
Number of lines x: 41 y: 6 z: 6
Mesh BBOX=[0,40,0,5,0,5]
```


<span id="page-9-0"></span>Table 2: Directive properties. \* indicates a multiplicity that while syntactically valid will lead to a useless simulation!

```
Grid dimensions [cells]: 58 x 7 x 7
Grid size: 1400 cells
Inner grid: BBOX=[9,49,1,6,1,6]
Outer grid: BBOX=[1,57,1,6,1,6]
Ghost grid: BBOX=[0,58,0,7,0,7]
Minimum edge lengths: DXMIN = 9.999983e-04, DYMIN = 9.999999e-04, DZMIN = 9.999999e-04
Maximum edge lengths: DXMAX = 1.000002e-03, DYMAX = 1.000000e-03, DZMAX = 1.000000e-03
Time step [s]: 1.667820e-12
PML XLO Region: BBOX=[1,8,1,5,1,5]
PML XHI Region: BBOX=[49,56,1,5,1,5]
PML YLO Region: BBOX=[9,48,1,0,1,5]
PML YHI Region: BBOX=[9,48,6,5,1,5]
PML ZLO Region: BBOX=[9,48,1,5,1,0]
PML ZHI Region: BBOX=[9,48,1,5,6,5]
Total array allocation 258.5 kiB
```
This part of the log file contains the time-step and total array allocation. An example of the detailed memory usage report which follows the grid characteristics is:

Memory usage:





# <span id="page-10-0"></span>2 The computational mesh

# <span id="page-10-1"></span>2.1 Mesh version and title: VM, CE, GS

The version of the mesh format must be specified by the first directive in the mesh file using the VM directive:

```
VM <i: majorVersion>.<i: minorVersion>.<i: patchVersion>
```
Here the three integers define the major version number, minor version number and patch level of the mesh format. For mesh files conforming to the format specified by this manual the version numbers should match those on the front title page.

An optional single line description or title for the mesh can be defined after the VM directive using the CE directive:

#### CE <s: comment>

Section 1 of the mesh file is then closed using a GS directive:

GS

# <span id="page-10-2"></span>2.2 Mesh lines and cells: DM, MS, XL, YL, ZL



<span id="page-10-3"></span>

The structured mesh used by Vulture is defined by a set of mesh lines  $x_i$ ,  $y_j$  and  $z_k$  in the x, y and z directions

$$
\begin{aligned}\n\{x_i; \quad i = 0 \dots N_x - 1\} \\
\{y_j; \quad j = 0 \dots N_y - 1\} \\
\{z_k; \quad k = 0 \dots N_z - 1\},\n\end{aligned} \tag{2.1}
$$

where  $N_x$ ,  $N_y$  and  $N_z$  are the number of mesh lines in each direction (see Figure [1\)](#page-10-3). The mesh lines are labelled by the indices  $i, j$  and  $k$ . The intersection points of the mesh lines define a set of nodes labelled by the indices of the intersecting mesh lines,  $(i, j, k)$ , and the lines (edges) between the nodes partition the volume of the mesh into  $(N_x - 1) \times (N_y - 1) \times (N_z - 1) \equiv n_x \times n_y \times n_z$  cuboid cells. These cell form the primary grid or lattice. The cells are labelled by the indices of the mesh lines on the low coordinate side of the cell, so we have cells number indices  $0, \ldots, n_x - 1$  etc. The cell sizes are given by

$$
\{\Delta x_i = x_{i+1} - x_i; i = 0 \dots n_x - 1\}
$$
  

$$
\{\Delta y_j = y_{j+1} - y_j; j = 0 \dots n_y - 1\}.
$$
  

$$
\{\Delta z_k = z_{k+1} - z_k; k = 0 \dots n_z - 1\}
$$
  
(2.2)

The electric field is sampled at the centres of the primary grid edges at times  $t = n\Delta t$  ( $n = 0, \ldots, N_s - 1$ ):  $E_x^n(i+\frac{1}{2},j,k)$ ,  $E_y^n(i,j+\frac{1}{2},k)$  and  $E_z^n(i,j,k+\frac{1}{2})$ . The cell centres are at the coordinates

$$
\begin{aligned}\n\{ \quad & x_{i+\frac{1}{2}} = x_i + \Delta x_i/2; \quad i = 0 \dots N_x - 2 \\
\{ \quad & y_{j+\frac{1}{2}} = y_j + \Delta y_j/2; \quad j = 0 \dots N_y - 2 \\
\{ \quad & z_{k+\frac{1}{2}} = z_k + \Delta z_k/2; \quad k = 0 \dots N_z - 2\n\}.\n\end{aligned}\n\tag{2.3}
$$

The *dual* or *secondary lattice* is then defined to be the grid whose vertices are the primary lattice centres,

{ 
$$
x_{i+\frac{1}{2}};
$$
  $i = 0...N_x - 2$ }  
\n{  $y_{j+\frac{1}{2}};$   $j = 0...N_y - 2$ }  
\n{  $z_{k+\frac{1}{2}};$   $k = 0...N_z - 2$ }, (2.4)

and the edge lengths of the dual lattice are

$$
\begin{aligned}\n\{\n\quad \tilde{\Delta}x_i &= x_{i+\frac{1}{2}} - x_{i-\frac{1}{2}} = (\Delta x_i + \Delta x_{i-1})/2; \quad i = 1 \dots N_x - 2 \\
\{\n\quad \tilde{\Delta}y_j &= y_{j+\frac{1}{2}} - y_{j-\frac{1}{2}} = (\Delta y_j + \Delta y_{j-1})/2; \quad j = 1 \dots N_y - 2 \\
\{\n\quad \tilde{\Delta}z_k &= z_{k+\frac{1}{2}} - z_{k-\frac{1}{2}} = (\Delta z_k + \Delta z_{k-1})/2; \quad k = 1 \dots N_z - 2.\n\end{aligned}\n\tag{2.5}
$$

The magnetic field is sampled at the centres of the secondary grid edges at times  $t = (n + \frac{1}{2})\Delta t$  ( $n =$  $(0,\ldots,N_{\rm s}-1)$ :  $H_{x}^{n+\frac{1}{2}}(i,j+\frac{1}{2},k+\frac{1}{2})$ ,  $H_{y}^{n+\frac{1}{2}}(i+\frac{1}{2},j,k+\frac{1}{2})$  and  $H_{z}^{n+\frac{1}{2}}(i+\frac{1}{2},j+\frac{1}{2},k)$ . The locations of the electric and magnetic fields in the primary are is shown in Figure [2.](#page-12-1)

In the mesh file the overall mesh extents are specified in Section 1 using the numbers of cells  $(n_x, n_y, n_z)$ with the DM directive:

DM <i: nx> <i: ny> <i: nz>

The mesh sizes are defined in Section 3 of the mesh file. Uniform meshes are defined using an MS directive to specify the constant edge lengths  $\langle \mathbf{r}: \mathbf{delay} \rangle (\Delta x)$ ,  $\langle \mathbf{r}: \mathbf{delay} \rangle (\Delta y)$  and  $\langle \mathbf{r}: \mathbf{delay} \rangle (\Delta z)$  in the three coordinate directions:

MS <r: delx> [ <r: dely> <r: delz> ] ]

The edge lengths are specified in metres. If  $\langle r: \text{ delay} \rangle$  and/or  $\langle r: \text{ delay} \rangle$  are missing the edge length from the previous coordinate direction is used. A uniform cubic mesh can therefore be defined using

MS <r: del>

in which case  $\Delta x_i = \Delta y_i = \Delta z_k = \Delta (\forall i, j, k)$ . The physical coordinates of the mesh lines are defined automatically for a uniform mesh starting at  $x_0 = y_0 = z_0 = 0$  m. For non-uniform meshes the MS directive is replaced by XL, YL and ZL directives which explicitly give the physical coordinates of all the mesh lines in metres:



<span id="page-12-1"></span>Figure 2: Location of the field sampling points in the primary FDTD grid.

```
XL
<r: x[0]>
.....
\langle r: x[nx] \rangleYL
<r: y[0]>
.....
<r: y[ny]>
ZL
<r: z[0]>
.....
\langle r: z[nz] \rangle
```
# <span id="page-12-0"></span>2.3 Bounding boxes

Objects on the mesh are specified by a bounding box selector which references a physical model or other entity such as a source or observer. An inclusive bounding box consists of a set of six integer mesh line indices  $[i_{\text{lo}}, i_{\text{hi}}, j_{\text{lo}}, j_{\text{hi}}, k_{\text{lo}}, k_{\text{hi}}]$  that selects all the compatible mesh elements within and on its surface:

 $[i_{\text{lo}}, i_{\text{hi}}, j_{\text{lo}}, j_{\text{hi}}, k_{\text{lo}}, k_{\text{hi}}] = \{i : i_{\text{lo}} \le i \le i_{\text{hi}}, j : j_{\text{lo}} \le j \le j_{\text{hi}}, k : k_{\text{lo}} \le k \le k_{\text{hi}}\}.$ 

Such bounding boxes are able to select single nodes, edges, faces, surfaces and volumes on the mesh. Bounding boxes are required to be normal in the sense defined by

$$
i_{\rm lo} \le i_{\rm hi}
$$
  
\n
$$
j_{\rm lo} \le j_{\rm hi}
$$
  
\n
$$
k_{\rm lo} \le k_{\rm hi},
$$
  
\n(2.6)

and use the mesh line numbers from  $(0, 0, 0)$  to  $(n_x, n_y, n_z)$ . Some examples are:

2 2 3 3 4 4 # A node at (2,3,4). 2 2 3 3 4 5 # A z-directed edge. 2 2 3 3 4 7 # A z-directed line 3 edges long. 2 2 1 2 1 2 # An x-normal face. 2 2 1 5 1 5 # An x-normal 4x4 face surface. 2 3 3 4 4 5 # A single cell. 2 4 3 5 4 6 # A 2x2x2 cell volume.

### <span id="page-13-0"></span>2.4 Simulation time-step control: NT, CN

The number of time-step iterations to be performed,  $N_s$ , is specified by the  $\leq$  i: numSteps> parameter in the NT directive:

#### NT <i: numSteps>

The FDTD simulation will run for  $N_s$  iterations  $n = 0, \ldots, N_s - 1$  corresponding to times  $t = 0, \ldots, (N_s - 1)$  $1)\Delta t$ . The time-step is determined automatically from the mesh-lines using

$$
\Delta t = C_N \frac{1}{c_0 \sqrt{\left(\frac{1}{\Delta x_{\min}}\right)^2 + \left(\frac{1}{\Delta y_{\min}}\right)^2 + \left(\frac{1}{\Delta z_{\min}}\right)^2}}
$$
(2.7)

where  $\Delta x_{\min} = \min_i [\Delta x_i, \tilde{\Delta} x_i]$  etc. are the edge lengths of the smallest cell in the mesh. Thus the cell with the largest value of  $1/(\Delta x_{\min})^2 + 1/(\Delta y_{\min})^2 + 1/(\Delta z_{\min})^2$  in the mesh determines the time-step. Here the Courant-Fredrichs-Lewy Number (CFLN),  $C_N$ , must be in the range zero to one,  $0 < C_N \le 1$ , for the simulation to be stable. By default  $C_N = \sqrt{3}/2$ , in which case on a cubic mesh of edge length  $\Delta$  the time-step becomes

$$
\Delta t = \frac{\Delta}{2c_0} \tag{2.8}
$$

and wavefronts propagating in free-space normal to mesh faces advance by one cell in two time-steps.

The CFLN number can changed using the CN directive:

### CN <r: courantNumber>

Dispersion in the FDTD grid becomes greater as the CFLN is reduced. However, as the CFLN approaches unity the algorithm has more chance of becoming unstable, depending on which features of the code are being used.

# <span id="page-13-1"></span>3 Physical models

# <span id="page-13-2"></span>3.1 Physical models and bounding boxes as selectors

Physical models for volumetric, surface and line type materials are defined separately using MT, BT and WT directives respectively:

MT <t: name> <t: type> <param1> ... <paramN> BT <t: name> <t: type> <param1> ... <paramN> WT <t: name> <t: type> <param1> ... <paramN>

The parameters for these directives are generally independent of the mesh and specific to the physical model. The name-spaces of the tags for the voulme, surface and wire material types are distinct.

Three corresponding types of selector are defined for physical models on the mesh using MB, TB and TW directives:



These select volumes (blocks), surfaces (boundaries) and lines (wires) respectively. These selectors are used to define the mesh elements on which the physical models referenced by their tags are to be applied. The parameters are used to define how the model is applied within the bounding box.

Irrespective of the order in which the physical model directives are declared in Section 2 of the mesh file they are applied to the computational grid in the following order:

1. All MBs in the order found in the mesh.

2. All TBs in the order found in the mesh.

3. All TWs in the order found in the mesh.

4. All EXs in the order found in the mesh.

5. All PWs in the order found in the mesh.

As will be detailed below MB, TB and TW objects of type PEC, FREE SPACE and SIMPLE will over-write each other without side effects. Other types (e.g SIBCs and dispersive materials) will not overwrite and will probably generate indeterminate results.

## <span id="page-14-0"></span>3.2 Volumetric materials: MT, MB

Volumetric material model types are defined using the MT directive which has the general form:

MT <t: name> <t: type> <param1> ... <paramN>

Here <t: name> is the name tag for the material type being defined and must be unique amongst all other MT directives. The valid material types,  $\lt t$ : type>, and their parameters are:

```
MT <t: name> SIMPLE [ <r: eps_r> [ <r: sigma> [ <r: mu_r> ] ] ]
MT <t: name> FREE SPACE
MT <t: name> PEC
MT <t: name> DEBYE <r: eps_Inf> <r: sigma> <r: mu_r> <r: res_1> <r: pole_1> \
   [ <r: res_2> <r: pole_2> [ <r: res_3> <r: pole_3> ] ]
MT <t: name> DEBYE <s: fileName>
```
The material type tags PEC and FREE SPACE are predefined for PEC and FREE SPACE materials respectively and cannot be overridden, i.e. the directives

MT PEC PEC MT FREE SPACE FREE SPACE

have effectively already been applied internally by the solver.

A material type identified by tag  $\leq t$ : name> is applied to a volume on the mesh using the MB directive:

MB <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> [ <t: mask> ]

The bounding box must define a volume (not a surface, line, edge or node). Volumetric materials are the first physical models applied to computational grid with each MB directive applied in the order it appears in the mesh file. Materials of type SIMPLE, FREE SPACE and PEC will overwrite each other without adverse side effects. Other materials types should not be applied "on top of each other". If compilation option USE AVERAGED MEDIA=OFF (see below) then the optional parameter <t: mask> can be used to determine if each external surface of the block is included in the material. The parameter is a string of six binary digits 0/1 with no spaces between them indicating whether (1) or not (0) the surfaces XLO, XHI, YLO, YHI, ZLO, ZHI respectively of the block are within the material. The default is to include all surfaces, i.e. 111111. The string 111011 would indicate that all surfaces except YHI should be included.

Depending on the value of the compilation option USE AVERAGED MEDIA, volumetric materials are applied to the grid in one of two ways:

- USE AVERAGED MEDIA=OFF: The simple material parameters are applied directly to all the edges encompassed by the bounding box. This method does not accurately represent the location of interfaces between different media. SIMPLE material types will overwrite when the bounding boxes overlap. Overlapping of DEBYE material types is ill-defined and the  $\leq t$ : mask> option of the MB directive should be used to align such material boundaries precisely without overlap.
- USE AVERAGED MEDIA=ON: The material parameters are first associated with the entire volume of each cell in the bounding box, ignoring any  $\leq t$ : mask> parameters in MB directives. After all the MB directives have been processed the parameters are then averaged across the boundaries of different materials to give a more accurate model for material interfaces at the location of the mesh faces (i.e the actual location of the bounding box surfaces). This option does not yet correctly support averaging of DEBYE type media. For Debye media the high frequency permittivity and (ionic) conductivity terms are averaged, however, the polarisation currents due to the relaxtion terms are not. The polarisation currents are applied directly, as for the case USE AVERAGED MEDIA=OFF, respecting any <t: mask> parameters. Depending on the material parameters and simulation geometry this may or may not produce an accurrate model.

# <span id="page-15-0"></span>3.2.1 SIMPLE non-dispersive material type

```
MT <t: name> SIMPLE [ <r: eps_r> [ <r: sigma> [ <r: mu_r> ] ] ]
```


<span id="page-15-3"></span>Table 3: SIMPLE material model parameters.

The simple material model type, SIMPLE, defines frequency independent isotropic media that are characterised by a relative permittivity,  $\epsilon_r$ , electrical conductivity,  $\sigma$ , and a relative permeability,  $\mu_r$ . The parameters for this type are defined in Table [3.](#page-15-3)

An example of a simple material model type and its application on a mesh is:

MT copper SIMPLE 1.0 3.7e7 1.0 MB 10 20 10 20 10 20 copper

This would set the volume enclosed by the bounding box [10, 20, 10, 20, 10, 20] to have the simple material parameters of copper:  $\epsilon_r = 1$ ,  $\sigma = 37$  MS/m and  $\mu_r = 1$ .

### <span id="page-15-1"></span>3.2.2 FREE SPACE material type

MT <t: name> FREE\_SPACE

The free-space material type, FREE SPACE, is a pseudonym for SIMPLE 1.0 0.0 1.0 so the above is equivalent to

MT <t: name> SIMPLE 1.0 0.0 1.0

### <span id="page-15-2"></span>3.2.3 PEC material type

MT <t: name> PEC

The PEC material type, PEC, simulates a perfect electric conductor in volumes to which it is applied by setting the FDTD update coefficients to  $\alpha = -1.0$  and  $\beta = 0.0$  (see the Code Implementation Manual for details [\[4\]](#page-66-1)). This method is more robust than using a high conductivity to represent a PEC material as in

MT POORPEC SIMPLE 1.0 1e5

but it behaves in the same way with regard to overwriting of overlapping bounding boxes and the averaging of neighbouring materials discussed above.

#### <span id="page-16-0"></span>3.2.4 DEBYE electrically dispersive material type

```
MT lt: name> DEBYE lt:r: eps_Inf> lt:r: sigma> lt:r: mu_r> lt:r: res_1> lt:r: pole_1> \
   [ <r: res_2> <r: pole_2> [ <r: res_3> <r: pole_3> ] ]
```

```
MT <t: name> DEBYE <s: fileName>
```
The DEBYE material type defines a frequency dependent electrically dispersive material characterised by a multi-pole generalised Debye dispersion relationship. The dispersion relationship for an  $N_m$ -pole relaxation is given by

$$
\hat{\epsilon}_r(\omega) = \epsilon_\infty - \frac{\jmath\sigma}{\omega \epsilon_0} + \sum_{k=1}^{N_m} \left\{ \frac{r_k}{\jmath\omega - p_k} + \frac{r_k^*}{\jmath\omega - p_k^*} \right\},\tag{3.1}
$$

where  $\epsilon_{\infty}$  is the relative permittivity in the high frequency limit,  $\sigma$  is an ionic conductivity term,  $r_k$  is the residue of the k-th pole and  $p_k$  is the k-th pole. Using this formulation the poles are enforced to occur in complex conjugate pairs. Such parameterisations can, for example, be obtained using a vector fitting algorithm [\[5\]](#page-66-2). A standard Debye relaxation is a special case with a real residue and pole given by

$$
p_k = -\frac{1}{\tau_k} \tag{3.2}
$$

$$
r_k = \frac{\Delta \epsilon_k}{2\tau_k} \tag{3.3}
$$

$$
\left\{\frac{r_k}{\jmath\omega - p_k} + \frac{r_k^*}{\jmath\omega - p_k^*}\right\} \rightarrow \frac{\Delta \epsilon_k}{1 + \jmath\omega \tau_k},\tag{3.4}
$$

where  $\Delta \epsilon_k = \epsilon_{\text{static};k} - \epsilon_{\infty;k}$  is the decrement caused by the k-th pole and  $\tau_k$  is the relaxtion time of the pole. Here  $\epsilon_{s,k}$  is the static, low frequency permittivity associated with the k-th relaxation. A second-order Lorentz dispersion is given by

$$
p_k = -\delta_k - \jmath \sqrt{\omega_k^2 - \delta_k^2} \tag{3.5}
$$

$$
r_k = \frac{\jmath \Delta \epsilon_k \omega_k^2}{2 \sqrt{\omega_k^2 - \delta_k^2}} \tag{3.6}
$$

$$
\left\{\frac{r_k}{\jmath\omega - p_k} + \frac{r_k^*}{\jmath\omega - p_k^*}\right\} \rightarrow \frac{\Delta \epsilon_k \omega_k^2}{\omega_k^2 + 2\jmath\omega \delta_k - \omega^2}.
$$
\n(3.7)

The parameters are summarised in Table [4,](#page-16-2) where the poles and residues are taken to be real. For a stable causal model it is required that  $\epsilon_{\infty} \geq 1$  and  $\text{Re}(p_k) \leq 0$ . For up to three real poles the Debye parameters can be specified directly in the mesh file. For complex poles and higher orders the parameters must be specified in an external ASCII file, referenced by the file name parameter  $\leq s$ : fileName>, that has the format given in Table [5.](#page-17-1) Note that only one residue/pole of each complex conjugate pair should be entered - the other is included implicitly!

| Parameter                         | Quantity                             | Symbol              | Units        | Default |
|-----------------------------------|--------------------------------------|---------------------|--------------|---------|
| $\langle r: \text{eps} \rangle$   | High frequency relative permittivity | $\epsilon_{\infty}$ |              | $1.0\,$ |
| $\langle r: \text{sigma} \rangle$ | (Ionic) electrical conductivity      | $\sigma$            | S/m          | 0.0     |
| $\langle r: res_k \rangle$        | Residue                              | $r_k$               | rad $s^{-1}$ | none    |
| $\langle r: pole_k\rangle$        | Pole                                 | $p_k$               | rad $s^{-1}$ | none    |

<span id="page-16-2"></span>Table 4: DEBYE material model parameters.

### <span id="page-16-1"></span>3.3 Internal and external surface materials, BT, TB

Surface material model types - called boundary types - are defined using the BT directive which has the general form:

### BT <t: name> <t: type> <param1> ... <paramN>

Here  $\lt t$ : name> is the name tag for the boundary type being defined and must be unique amongst all other BT directives. The valid boundary types, <t: type>, and their parameters are:

```
<i: N_m> <r: eps_inf> <r: sigma> <r: mu_r>
Re[\langle r: res\_1 \rangle] Im[\langle r: res\_1 \rangle] Re[\langle r: pole\_1 \rangle] Im[\langle r: pole\_1 \rangle]... ... ... ...
Re[<r: res_N>] Im[<r: res_N>] Re[<r: pole_N>] Im[<r: pole_N>]
```
<span id="page-17-1"></span>Table 5: External file format for the DEBYE material type. Comment and blank lines are not allowed.

BT <t: name> FREE\_SPACE BT <t: name> PEC BT <t: name> PMC BT <t: name> PERIODIC BT <t: name> MUR BT <t: name> PML <i: nlayer> <i: order> <rl: n\_eff> <rl: refcoeff> <rl: kmax> BT  $\lt t$ : name> SIBC  $\lt r$ : S11\_TM>  $\lt r$ : S12\_TM>  $\lt r$ : S21\_TM>  $\lt r$ : S22\_TM>  $\setminus$ [ <r: S11\_TE> <r: S12\_TE> <r: S21\_TE> <r: S22\_TE> ] BT <t: name> SIBC <s: fileName>

The boundary type tags PEC, PMC and FREE SPACE are predefined for PEC, PMC and FREE SPACE boundary types respectively and cannot be overridden, i.e. the directives

```
BT PEC PEC
BT PMC PMC
BT FREE_SPACE FREE_SPACE
```
are effectively applied internally by the solver. The valid usage of the different boundary types on internal and external surfaces of the mesh is summarised in Table [6.](#page-17-2)



<span id="page-17-2"></span>Table 6: Valid usage of different boundary types.

A boundary type identified by tag  $\lt t$ : name> is applied to a surface on the mesh using the TB directive:

```
TB \langle i: ilo \rangle \langle i: ihi \rangle \langle i: ilo \rangle \langle i: jhi \rangle \langle i: khio \rangle \langle i: khi \rangle \langle t: name \rangle[ <i: orient> [ <r: angle> ] ]
```
The bounding box must define a surface (not a volume, line, edge or node). The parameters  $\leq i$ : orient> and  $\langle r: \text{ angle}\rangle$  are used to orientate and align asymmetric and anisotropic boundary type materials on the computational grid as discussed below.

Boundary type materials are the second type of physical model applied to computational grid, after the volumetric materials, with each TB directive applied in the order it appears in the mesh file. Boundaries of type FREE SPACE and PEC will overwrite each other without adverse side effects. SIBC boundaries can be applied (once) on top of PEC boundaries. Other boundary types should not be applied "on top of each other".

# <span id="page-17-0"></span>3.3.1 External surfaces

Boundary type tags for the boundaries forming each external face of the mesh are predefined by the solver and associated with those surfaces. This is equivalent to the directives



having been applied, where  $nx$ ,  $ny$ ,  $nz$  are the mesh extents. So, for example, XLO is a tag for the boundary type of the x-normal external face of the mesh on the lower  $x$  side of the mesh. These tags can used to define the boundary types on the external mesh surfaces using BT directives, somewhat in reverse to the way they are used for all other internal surfaces. By default all the external surfaces are PML absorbing boundary types with default parameters, i.e. the directives

BT XLO PML BT XHI PML BT YLO PML BT YHI PML BT ZLO PML BT ZHI PML

are applied by default. Only boundary types PEC, PMC, PERIODIC, MUR and PML are valid on external surfaces.

## <span id="page-18-0"></span>3.3.2 PEC boundary type

#### BT <t: name> PEC

The PEC boundary type enforces zero tangential electric field on a surface by setting the FDTD update coefficients to  $\alpha = -1.0$  and  $\beta = 0.0$  (see the Code Implementation Manual for details [\[4\]](#page-66-1)).

The following example defines a closed PEC box:

BT walls PEC TB 10 10 10 20 10 20 walls TB 20 20 10 20 10 20 walls TB 10 20 10 20 10 10 walls TB 10 20 10 20 20 20 walls TB 10 20 10 10 10 20 walls TB 10 20 20 20 10 20 walls

# <span id="page-18-1"></span>3.3.3 PMC boundary type

#### BT <t: name> PMC

The PMC boundary type enforces a PMC boundary condition on an external surface using a field mirroring algorithm with ghost fields "outside" the computational grid. In this way the PMC is located on the primary grid faces, the same as PEC boundary types on external surfaces.

For example an infinite parallel plate waveguide with metal plates on the lower and upper z-normal external surfaces can be defined using:

BT XLO PMC BT XHI PMC BT YLO PEC BT YHI PEC BT ZLO PML BT ZHI PML

PMC boundary types can also be used to create symmetry planes, for example:

BT XLO PML BT XHI PML BT YLO PMC BT YHI PML BT ZLO PML BT ZHI PML

would create a symmetry plane in the YLO mesh surface.

# <span id="page-19-0"></span>3.3.4 PERIODIC boundary type

#### BT <t: name> PERIODIC

The PERIODIC boundary type enforces a periodic boundary condition between pairs of external surfaces XLO/XHI, YLO/YHI and ZLO/ZHI, using a field copying algorithm with ghost fields "outside" the computational grid. The boundary condition is located on the primary grid faces, the same as other boundary types on external surfaces. Opposite pairs of external faces must both be defined with the PERIODIC boundary type.

For example an infinite parallel plate waveguide with metal plates on the lower and upper z-normal external surfaces can be defined using:

BT XLO PERIODIC BT XHI PERIODIC BT YLO PEC BT YHI PEC BT ZLO PML BT ZHI PML

### <span id="page-19-1"></span>3.3.5 FREE SPACE boundary type

BT <t: name> FREE\_SPACE

The FREE SPACE boundary type imposes free-space material parameters on a surface. The main use of this boundary type is to "punch" holes through internal PEC type surfaces to create apertures. For example:

```
BT walls PEC
BT aperture FREE_SPACE
# First make a closed metal box.
TB 10 10 10 20 10 20 walls
TB 20 20 10 20 10 20 walls
TB 10 20 10 20 10 10 walls
TB 10 20 10 20 20 20 walls
TB 10 20 10 10 10 20 walls
TB 10 20 20 20 10 20 walls
# Then punch hole in face.
TB 10 10 12 18 12 18 aperture
```
Note that this does not work for SIBC types surfaces!

#### <span id="page-19-2"></span>3.3.6 MUR boundary type

#### BT <t: name> MUR

The MUR boundary type specifies a first order Mur analytic ABC for external surfaces. It is compatible with PEC and PMC boundary types on neighbouring surfaces. It gives a normal incidence free-space reflection coefficient of approximately -20 dB providing sufficient space is left between it and scattering objects on the grid (about half a wavelength).

### <span id="page-19-3"></span>3.3.7 PML boundary type

#### BT <t: name> PML <i: nlayer> <i: order> <r: n\_eff> <r: refcoeff> <r: kmax>

The PML boundary type specifies a perfectly matched layer (PML) ABC for external surfaces. This is a more advanced ABC than the Mur with much better performance, albeit at a significantly higher computational cost. The PML boundary type adds a number of layers,  $n_{\text{layer}}$ , of fictitious matched absorbing material to the external face of the mesh. Adding more layers improves the performance, i.e. reduces the reflections from the face, but increases the memory requirements and run-time of the simulation. Vulture implements a Uniaxial PML (UPML) with complex stretching factor [\[4\]](#page-66-1)

$$
s(x) = \kappa(x) + \frac{\sigma(x)}{\mathrm{i}\omega\epsilon_0} \tag{3.8}
$$

and polynomial grading profiles

$$
\sigma(x) = \sigma^{\max}\left(\frac{x}{d}\right)^m \tag{3.9}
$$

$$
\kappa(x) = 1 + (\kappa^{\max} - 1) \left(\frac{x}{d}\right)^m.
$$
\n(3.10)

Here  $x$  is the perpendicular distance into the PML and  $d$  is the total thickness of the PML. The maximum PML conductivity,  $\sigma^{\text{max}}$ , is related to the magnitude of the theoretical (continuous space) reflection coefficient,  $|R(0)|$  by

$$
\sigma^{\max} = -\frac{m+1}{2\eta d} \ln |R(0)|. \tag{3.11}
$$

This has an empirically determined optimum value of

$$
\sigma^{\text{opt}} = \frac{4}{\eta \Delta} \frac{m+1}{5} = \frac{4n_{\text{eff}}}{\eta_0 \Delta l} \frac{m+1}{5},\tag{3.12}
$$

where  $\Delta$  is the mesh size of the PML cells in the direction perpendicular to the surface and  $n_{\text{eff}}$  is the effective refractive index of the dominant mode(s) incident on the PML. The PML parameters and their default values are defined in Table [7.](#page-20-1)

| Parameter                           | Quantity                   | Symbol             | <b>Units</b> | Default           |
|-------------------------------------|----------------------------|--------------------|--------------|-------------------|
| $\langle i: \text{ nlayer} \rangle$ | Number of layers           | $n_{\rm layer}$    |              |                   |
| $\langle i: order \rangle$          | Profile order              | m                  |              |                   |
| $\langle r: n_{eff} \rangle$        | Effective refractive index | $n_{\text{eff}}$   |              | 1.0               |
| $\langle r:$ refcoeff>              | Reflection coefficient     | R(0)               |              | $-1$ <sup>*</sup> |
| $\langle r: \text{kmax}\rangle$     | Maximum PML permittivity   | $\kappa_{\rm max}$ |              | $1.0\,$           |

<span id="page-20-1"></span>Table 7: PML boundary type model parameters.  $*$  indicates optimal value should be chosen automatically

While the PML has substantially improved performance relative to the Mur ABC it has limitiations:

- The reflection coefficient begins to increase at grazing angles of incidence which can be a problem on high apsect ratio meshes.
- The PMLs performance at transparently terminating evanescent waves is limited. The PML permittivity parameter  $\kappa$  is useful for improving the performance in this respect. Typically values of  $\kappa$  from 5 to 80 have been found to be useful.

The Vulture PML implementation allows inhomogeneous simple materials to be place in contact with PML boundaries and PEC type internal boundaries normal to the PML surface can also be placed in contact with the PML. This effectively simulates the medium touching the PML being extended in the normal direction to infinity. The PML is also consistent with PEC, PMC and PERIODIC boundary types on neighbouring external surfaces. Support for electrically dispersive materials is currently only partially implemented.

For example, a 10 layer PML with a cubic polynomial profile can de defined using:

BT ZLO PML 10 3

<span id="page-20-0"></span>3.3.8 SIBC boundary type

```
BT <t: name> SIBC <s: fileName>
BT ltt: name> SIBC ltr: Saa_TM> ltr: Sab_TM> ltr: Sba_TM> ltr: Sbb_TM> \setminus[ <r: Saa_TE> <r: Sab_TE> <r: Sba_TE> <r: Sbb_TE> ]
TB \langle i: ilo \rangle \langle i: ihi \rangle \langle i: jlo \rangle \langle i: jhi \rangle \langle i: klo \rangle \langle i: khi \rangle \langle t: name \rangle[ <i: orient> [ <r: angle> ] ]
```
Note that the SIBC boundary type is a compile time option that is enabled by setting the compilation option WITH SIBC=ON and may not be available in a particular solver executable.

SIBC boundary types are arbitrary frequency dependent reflection and transmission boundaries. In general the tangential fields on each side of an anisotropic boundary representing a thin material can be



<span id="page-21-0"></span>Figure 3: Definition of SIBC impedance matrix (left) and alignment SIBCs on the mesh for a z-normal surface(right). The conventions for  $x$ - and  $y$ -normal surface are given by cyclic permutation.

resolved into transverse electric (TE) and transverse magnetic (TM) waves (see Figure [3\)](#page-21-0). The SIBC model relates these tangential electric and magnetic fields using a surface impedance matrix. A natural convention for defining the surface impedance matrix is with the TM and TE ports blocked together on either side of the boundary. In the principal axis system of the material the impedance matrix is block diagonal and the SIBC can be written

$$
\begin{bmatrix} E_a^{\text{TM}}(\omega) \\ E_b^{\text{TM}}(\omega) \\ E_b^{\text{TE}}(\omega) \\ E_b^{\text{TE}}(\omega) \end{bmatrix} = \begin{bmatrix} Z_{ab}^{\text{TM}}(\omega) & Z_{ab}^{\text{TM}}(\omega) & 0 & 0 \\ Z_{ab}^{\text{TM}}(\omega) & Z_{bb}^{\text{TM}}(\omega) & 0 & 0 \\ 0 & 0 & Z_{bb}^{\text{TE}}(\omega) & Z_{ab}^{\text{TE}}(\omega) \\ 0 & 0 & Z_{ba}^{\text{TE}}(\omega) & Z_{bb}^{\text{TE}}(\omega) \end{bmatrix} \begin{bmatrix} H_a^{\text{TM}}(\omega) \\ -H_b^{\text{TM}}(\omega) \\ H_a^{\text{TM}}(\omega) \\ H_a^{\text{TE}}(\omega) \\ -H_b^{\text{TE}}(\omega) \end{bmatrix} \equiv Z(\omega) \begin{bmatrix} H_a^{\text{TM}}(\omega) \\ -H_b^{\text{TM}}(\omega) \\ H_a^{\text{TE}}(\omega) \\ -H_b^{\text{TE}}(\omega) \end{bmatrix} (3.13)
$$

with two 2-by-2 sub-matrices for the TM and TE polarisations. Here "a" and "b" denote the two sides of the boundary. For a reciprocal material

$$
Z_{ij}(\omega) = Z_{ji}(\omega),\tag{3.14}
$$

while if the material is isotropic

$$
Z_{ij}^{\text{TM}}(\omega) = Z_{ij}^{\text{TE}}(\omega) \equiv Z_{ij}(\omega). \tag{3.15}
$$

The TM and TE polarisations are defined with reference to a local reference system on the surface of the material. Aligning this with the global coordinate system of a computational grid requires two parameters to be specified for each surface on which the SIBC is applied:

- $\leq$  i: orient> ( $\xi$ ): A binary orientation flag that indicates which side of the material is on which side of the surface.  $\xi = +1$  indicates the "a" side of the material is on the lower perpendicular coordinate side of the surface while  $\xi = -1$  indicates the converse.
- $\langle \mathbf{r}: \mathbf{angle} \rangle$  ( $\alpha$ ): The angle between the  $x/y/z$  coordinate axis on the mesh and the  $x_{pa}$  reference axes on the material for  $z/x/y$ -normal surfaces respectively.

The conventions are shown in Figure [3](#page-21-0) and the SIBC parameters and their default values are summarised in Table [8.](#page-22-0) SIBC boundary types can only be used on internal surfaces.

The impedance matrix is related to the scattering matrix of the surface by

$$
S = (Z - Z_0)(Z + Z_0)^{-1}, \tag{3.16}
$$

where

$$
Z_0 = \text{diag}\left[\eta_{\text{TM}}^a, \eta_{\text{TE}}^b, \eta_{\text{TE}}^a, \eta_{\text{TE}}^b\right]
$$
\n(3.17)

| Parameter                         | Quantity                       | Symbol   | Units   | Default |
|-----------------------------------|--------------------------------|----------|---------|---------|
| $\leq$ : filename>                | File name                      | n/a      |         |         |
| <i: orient=""></i:>               | Orientation of sides $(\pm 1)$ |          |         | $+1$    |
| $\langle r: \text{ angle}\rangle$ | Angle of principal axes        | $\alpha$ | degrees | 0.0     |

<span id="page-22-0"></span>Table 8: SIBC boundary type model parameters.

is a diagonal matrix holding the intrinsic impedances of the media on either side of the boundary. Note that the scattering matrix is specific to the medium surrounding the boundary whereas the impedance matrix is not.

The frequency dependent impedance matrix elements are represented by partial fraction expansions (PFEs) in the s-plane

$$
Z_{ij}(s = j\omega) = Z_{ij}^{\infty} + \sum_{k=1}^{N_{ij}} \frac{r_{ij}^k}{s - p_{ij}^k} \quad (i, j = 1, ..., M)
$$
\n(3.18)

where  $Z_{ij}^{\infty}$  is the high frequency asymptotic response,  $N_{ij}$  is the order of the  $(i, j)$ -th element and  $r_{ij}^k$  and  $p_{ij}^k$  are the k-th complex residues and poles of the  $(i, j)$ -th element. Typically the order is the same for each element, however, this does not need to be the case. Frequency independent boundaries are a special case with  $N_{ij} = 0 \, (\forall i, j)$ .

The PFE representation automatically ensures causality of the SIBC. Stability requires that the poles all lie in the left-hand side of the complex plane  $(\text{Re}[p_{ij}^k] < 0)$  and passivity requires that  $\text{Re}[Z_{ij}(\omega)] \geq 0 \ \forall \omega$ . The latter condition can be difficult to ensure since it applies to all frequencies, including those far above the frequency at which the mesh becomes under-sampled.

The PFEs are specified in an external ASCII format file referenced by the  $\leq$ s: filename> parameter. Denoting  $r_{ij}^k$  by  $\mathbf{r}(i,j,k)$ ,  $p_{ij}^k$  by  $\mathbf{p}(i,j,k)$  and  $N_{ij}$  by  $\mathbb{N}(i,j)$  the general format of this PFE input file is given in Table [9.](#page-23-1) Vulture currently supports three cases of SIBC in external files determined by the order of the impedance matrix, M:

- $M = 1$ : NOT IMPLEMENTED YET: One-sided (no transmission) isotropic symmetrical material.
- $M = 2$ : Two-sided isotropic material.
- $M = 4$ : Two-sided anisotropic material.

For frequency independent boundaries real valued scattering parameters

$$
\begin{bmatrix} E_{\rm a}^{\rm TM} \\ E_{\rm b}^{\rm TM} \\ E_{\rm a}^{\rm TE} \\ E_{\rm b}^{\rm TE} \end{bmatrix} = \begin{bmatrix} S_{aa}^{\rm TM} & S_{ab}^{\rm TM} & 0 & 0 \\ S_{ba}^{\rm TM} & S_{bb}^{\rm TM} & 0 & 0 \\ 0 & 0 & S_{ab}^{\rm TE} & S_{bb}^{\rm TE} \\ 0 & 0 & S_{bb}^{\rm TE} & S_{bb}^{\rm TE} \\ 0 & 0 & S_{bb}^{\rm TE} & S_{bb}^{\rm TE} \end{bmatrix} \begin{bmatrix} E_{\rm a}^{\rm TM} \\ E_{\rm b}^{\rm TM} \\ E_{\rm b}^{\rm TE} \\ E_{\rm a}^{\rm TE} \\ E_{\rm b}^{\rm TE} \end{bmatrix} \equiv \mathcal{S} \begin{bmatrix} E_{\rm a}^{\rm TM} \\ E_{\rm b}^{\rm TM} \\ E_{\rm b}^{\rm TE} \\ E_{\rm a}^{\rm TE} \\ E_{\rm b}^{\rm TE} \end{bmatrix} \tag{3.19}
$$

can be specified directly in the mesh file using

$$
\begin{array}{lcl} \texttt{BT < t: name>} \ \texttt{SIBC} < r: \ \texttt{Saa_TM> < r: \ Sab_TM> < r: \ Sba_TM> < r: \ Sbb\_TH> < \\ & \texttt{[} < r: \ \texttt{Saa_TE> < r: \ Sab_TE> < r: \ Sba_TE> < r: \ Sbb\_TE> < \\ \end{array}
$$

These scattering parameters should be referenced to free-space port impedances, even if they are to be used on surfaces embedded in other media. The scattering matrix must represent a passive system, i.e.  $I - S^{\dagger}S$ must be positive definite, where I is the identity matrix.

For example, a surface composed of a frequency independent, isotropic, symmetric and reciprocal surface material with a reflection coefficient of -6 dB and transmission coefficient -20 dB can be modelled using

BT absorber SIBC -0.5 0.1 0.1 -0.5 TB 10 10 12 18 12 18 absorber

An example of an enclosure with five metal walls and one wall composed of a frequency dependent surface material defined by the PFE in a file called "cfc.prm" is:



<span id="page-23-1"></span>Table 9: PFE ASCII file format for SIBC boundary types. For Vulture SIBC boundaries only  $M = N = 2, 4$ are currently valid.

BT cfc2 SIBC "cfc.prm" BT walls PEC TB 10 10 10 20 10 20 walls TB 20 20 10 20 10 20 walls TB 10 20 10 20 10 10 walls TB 10 20 10 20 20 20 walls TB 10 20 10 10 10 20 walls TB 10 20 20 20 10 20 walls TB 10 10 12 18 12 18 cfc2

# <span id="page-23-0"></span>3.4 Linear wire materials: WT, TW

Wire material types are defined using the  $WT$  directive which has the general form:

WT <t: name> <t: type> <param1> ... <paramN>

Here <t: name> is the name for the wire type being defined and must be unique amongst all other WT directives. The valid wire types and their parameters are:

WT <t: name> PEC WT <t: name> THIN\_WIRE <r: radius>

The wire type tag PEC is predefined for PEC wire types and cannot be overridden, i.e. the directive

WT PEC PEC

has effectively already been applied internally by the solver.

A wire type identified by tag  $\leq t$ : name> is applied to a line on the mesh using the TW directive:

TW <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> \ [ <t: end1Type>] [ <t: end2Type> ]

The bounding box must define a line (not a volume, surface, or node). The parameters  $\lt t$ : endType1> and <t: endType2> are used to set the end types for the low and high ends of wire on the computational grid as discussed below.

Wire type materials are the third type of physical model applied to computational grid, after the volumetric and surface materials, with each TW directive applied in the order it appears in the mesh file. Wire types should not be applied "on top of each other". Wires can be applied within SIMPLE volumetric materials and across PEC type surfaces.

#### <span id="page-24-0"></span>3.4.1 PEC wire type

WT <t: name> PEC

The PEC wire types enforces zero electric field on a line by setting the FDTD update coefficients to  $\alpha = -1.0$  and  $\beta = 0.0$  (see the Code Implementation Manual for details [\[4\]](#page-66-1)). This method is more robust than using a high conductivity to represent a PEC material. This produces a model of a wire with crosssectional area of roughly 3 % of the the area of the secondary grid face normal to the wire's edge. Example:

```
BT cable PEC
TB 10 10 10 20 10 10 cable
```
#### <span id="page-24-1"></span>3.4.2 THIN WIRE wire type

The THIN\_WIRE type is currently not implemented!

```
WT <t: name> THIN_WIRE <r: radius>
TW <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> \
   [ <t: end1Type>] [ <t: end2Type> ]
```
The THIN WIRE type is a thin wire sub-cell model in which the wire radius,  $\langle \mathbf{r}: \mathbf{radius} \rangle$ , must be much less than the cell size. The radius is specified in metres. The parameters <t: endType1> and <t: endType2> are used to set the end types for the low and high coordinate ends of the wire. The valid values are THRU (default), END and CORNER.

An example for a wire going around a corner is:

WT cable THIN\_WIRE 1e-3 TW 10 10 10 20 10 10 cable END CORNER TW 10 10 20 20 10 20 cable CORNER END

# <span id="page-24-2"></span>4 Waveforms and sources

#### <span id="page-24-3"></span>4.1 Waveforms: WF

Waveform types are used to define a nominally dimensionless time-series of values that can be used by different physical sources in a simulation. They are defined by the WF directive which has two slightly different forms depending on the waveform type:

```
WF ltt: name> ltt: type> [ ltr: size> [ ltr: delay> [ ltr: width> [ ltr:freq> ] ] ]
WF <t: name> EXTERNAL <s: fileName> [ <r: size> [ <r: delay> ] ]
```
Here  $\lt t$ : name> is the name tag for the waveform type being defined and must be unique amongst all other WT directives. The valid waveform types and their parameters are

```
WF <t: name> GAUSSIAN_PULSE [ <r: size> [ <r: delay> [ <r: width> [ <r:freq> ] ] ]
WF <t: name> NARROW_GAUSSIAN_PULSE [ <r: size> [ <r: delay> [ <r: width> [ <r:freq> ] ] ]
WF <t: name> DIFF_GAUSSIAN_PULSE [ <r: size> [ <r: delay> [ <r: width> [ <r:freq> ] ] ]
WF \lt t: name> RICKER_WAVELET [\lt r: size> [\lt r: delay> [\lt r: width> [\lt r: freq> ] ] ]
WF ltt: name> COMPACT_PULSE [ ltr: size> [ ltr: delay> [ ltr: width> [ ltr:freq> ] ] ]
WF ltt: name> DIFF_COMPACT_PULSE [ ltr: size> [ ltr: delay> [ ltr: width> [ ltr:freq> ] ] ]
WF ltt: name> RAMPED_SINSOID [ ltr: size> [ ltr: delay> [ ltr: width> [ ltr:freq> ] ] ]
WF <t: name> MOD_GAUSSIAN_PULSE [ <r: size> [ <r: delay> [ <r: width> [ <r:freq> ] ] ]
WF ltt: name> MOD_COMPACT_PULSE [ ltr: size> [ ltr: delay> [ ltr: width> [ ltr:freq> ] ] ]
WF <t: name> EXTERNAL <s: fileName> [ <r: size> [ <r: delay> ] ]
```
The common waveform type parameters are defined in Table [10.](#page-25-1)

| Parameter                       | Quantity             | Symbol      | Units        | Default |
|---------------------------------|----------------------|-------------|--------------|---------|
| $\langle r: size \rangle$       | Magnitude            |             |              | $1.0\,$ |
| $\langle r: delay \rangle$      | Delay                | $t_{\rm d}$ | S            | 0.0     |
| $\langle r: width \rangle$      | Pulse width          |             |              | $\ast$  |
| $\langle r: \text{freq}\rangle$ | Modulation frequency | / m         | $_{\rm{Hz}}$ | $\ast$  |

<span id="page-25-1"></span>Table 10: Waveform type common parameters. <sup>∗</sup> depends on type.

#### <span id="page-25-0"></span>4.1.1 Built-in waveform types

The built-in waveform types are defined and their default parameters given in Table [11.](#page-28-3) The equations for the waveforms are:

Gaussian pulse:

$$
\psi_{\rm GP}(t) = A e^{-(t-t_{\rm d})^2/2\sigma^2} \tag{4.1}
$$

Differentiated Gaussian pulse:

$$
\psi_{\rm DGP}(t) = \sigma \frac{d\psi_{\rm GP}}{dt} = -A(t - t_{\rm d})e^{-(t - t_{\rm d})^2/2\sigma^2}
$$
\n(4.2)

Ricker wavelet (double-differentiated Gaussian pulse):

$$
\psi_{\rm RW}(t) = \sigma \frac{d\psi_{\rm DGP}}{dt} = \sigma^2 \frac{d^2 \psi_{\rm GP}}{dt^2} = -A \left( 1 - \left\{ \frac{t - t_{\rm d}}{\sigma^2} \right\}^2 \right) e^{-(t - t_{\rm d})^2 / 2\sigma^2}, \tag{4.3}
$$

Compact pulse (Blackman-Harris pulse):

$$
\psi_{\rm CP}(t) = \begin{cases} \frac{A}{32} (10 - 15 \cos(\omega_0 t) + 6 \cos(2\omega_0 t) - \cos(3\omega_0 t)) & 0 \le t \le 2\sigma \\ 0 & t > 2\sigma \end{cases}
$$
(4.4)

$$
= \begin{cases} \sum_{m=0}^{3} a_m \cos(m\omega_0 t) & 0 \le t \le 2\sigma \\ 0 & t > 2\sigma \end{cases}
$$
(4.5)

with 
$$
\omega_0 = \frac{\pi}{\sigma}
$$
,  $a_0 = 10/32$ ,  $a_1 = -15/32$ ,  $a_2 = 6/32$ ,  $a_3 = -1/32$  (4.6)

Differentiated compact pulse:

$$
\psi_{\text{DCP}}(t) = \sigma \frac{\mathrm{d}\psi_{\text{CP}}}{\mathrm{d}t} = \begin{cases} -A\frac{\pi}{32} \left( 15\sin(\omega_0 t) + 12\sin(2\omega_0 t) + 3\sin(3\omega_0 t) \right) & 0 \le t \le 2\sigma \\ 0 & t > 2\sigma \end{cases} \tag{4.7}
$$

Ramped sinusoid:

$$
\psi_{\rm RS}(t) = \begin{cases} A \sum_{m=0}^{3} a_m \cos(m\omega_0 t) \sin(2\pi f_m t) & 0 \le t \le \sigma \\ A \sin(2\pi f_m t) & t > \sigma \end{cases}
$$
(4.8)

Modulated Gaussian pulse:

$$
\psi_{\text{MGP}}(t) = A e^{-(t - t_{\text{d}})^2 / 2\sigma^2} \sin(2\pi f_m t) \tag{4.9}
$$

Modulated compact pulse:

$$
\psi_{\text{MCP}}(t) = \begin{cases} A \sum_{m=0}^{3} a_m \cos(m\omega_0 t) \sin(2\pi f_m t) & 0 \le t \le 2\sigma_{\text{CP}} \\ 0 & t > 2\sigma_{\text{CP}} \end{cases}
$$
(4.10)

When choosing a waveform the following issues should be considered:

- Broad pulses generally have more energy at high frequencies and therefore exhibit more artefacts due to under-sampling. The default parameters limit this effect while still giving sufficient energy at frequencies up to where cell size is  $\lambda/5$  to avoid dynamic range issues. The NARROW GAUSSIAN PULSE is an except to this - it has significant energy beyond the Nyquist spatial sampling frequency.
- Differentiated pulses have no DC response. The low frequency response may be inaccurate.
- The modulated Gaussian and compact pulse have a band-pass response, hence the results are only meaningful over a limited frequencies range and may be unreliable at low and high frequencies.

The default pulses and differentiated pulses and their spectra are shown in Figures [4](#page-26-0) and [5.](#page-26-1) The modulated pulses and their spectra are shown in Figure [6](#page-27-0) and Figure [7.](#page-27-1)



<span id="page-26-0"></span>Figure 4: Default pulse waveforms and their derivatives.



<span id="page-26-1"></span>Figure 5: Spectra of the pulse waveforms and their derivatives.



<span id="page-27-0"></span>Figure 6: Default modulated waveforms.



<span id="page-27-1"></span>Figure 7: Spectra of the default modulated waveforms.

| Waveform type         | Definition             | Size, A                  | Delay, $t_{\rm d}$ | Width, $\sigma$      | Frequency, $f_{\rm m}$  |
|-----------------------|------------------------|--------------------------|--------------------|----------------------|-------------------------|
|                       |                        | $\overline{\phantom{a}}$ | $\mathbf{S}$       | $\mathbf{s}$         | Hz)                     |
| GAUSSIAN_PULSE        | $\psi_{\text{GP}}(t)$  | 1.0                      | $40\Delta t$       | $5\sqrt{2}\Delta t$  | n/a                     |
| NARROW_GAUSSIAN_PULSE | $\psi_{\text{GP}}(t)$  | 1.0                      | $12\Delta t$       | $2\Delta t$          | n/a                     |
| DIFF_GAUSSIAN_PULSE   | $\psi_{\text{DGP}}(t)$ | 1.0                      | $40\Delta t$       | $5\sqrt{2}\Delta t$  | n/a                     |
| RICKER_WAVELET        | $\psi_{\text{RW}}(t)$  | 1.0                      | $40\Delta t$       | $5\sqrt{2}\Delta t$  | n/a                     |
| COMPACT_PULSE         | $\psi_{\rm CP}(t)$     | 1.0                      | n/a                | $20\Delta t$         | n/a                     |
| DIFF_COMPACT_PULSE    | $\psi_{\text{DCP}}(t)$ | 1.0                      | n/a                | $20\Delta t$         | n/a                     |
| RAMPED_SINUSOID       | $\psi_{\rm RS}(t)$     | 1.0                      | n/a                | $20\Delta t$         | $\overline{20\Delta t}$ |
| MOD_GAUSSIAN_PULSE    | $\psi_{\text{MGP}}(t)$ | 1.0                      | $120\Delta t$      | $20\sqrt{2}\Delta t$ | $20\Delta t$            |
| MOD_COMPACT_PULSE     | $\psi_{\rm MCP}(t)$    | 1.0                      | n/a                | $80\Delta t$         | $\overline{20\Delta t}$ |

<span id="page-28-3"></span>Table 11: Waveform types and default parameters.

### <span id="page-28-0"></span>4.1.2 EXTERNAL waveform type

A user defined waveform can be provided as a time-series in an external ASCII file using the EXTERNAL waveform type:

```
WF <t: name> EXTERNAL <s: fileName> [ <r: size> [ <r: delay> ] ]
```
<s: fileName> is the path to an ASCII file containing two columns with monotonically increasing time (seconds) and waveform values (-). The file format does not support any comment lines or blank lines! The waveform is assumed zero outside the range of times in the file and is interpolated using cubic splines for values in between those given in the file. The waveform is scaled by  $\langle r: size \rangle$  and delayed by  $\langle r: delay \rangle$ seconds.

# <span id="page-28-1"></span>4.2 Current and field sources: EX

Current and field sources are applied to the grid using the EX directive, which has two forms:

```
EX <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: type> \
   <t: wfName> [ <r: size> [ <r: delay> ] ]
EX <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: type> \
   <t: wfName> <r: param> [ <r: size> [ <r: delay> ] ]
```
The first form is valid for all the distributed field and current sources while the second form is used for lumped sources with finite internal impedance. Field and current sources are applied to the grid in the order they are defined in the mesh file before any plane-wave sources.

The common source parameters are defined in Table [12.](#page-28-4) The waveform to be used for the source, <t: wfName>, must have been previously defined in a WF directive. The delay of the source relative to its waveform,  $\langle r: \text{ delay} \rangle$ , is in addition to any delay specified in the WF directive.

| Parameter                 | Quantity      | Symbol         | Units  | Default |
|---------------------------|---------------|----------------|--------|---------|
| $lt:$ name>               | Source name   | n/a            | n/a    | none    |
| $< t$ : type $>$          | Source type   | n/a            | n/a    | none    |
| $lt:$ wfName>             | Waveform name | n/a            | n/a    | $\ast$  |
| $\langle r: size \rangle$ | Magnitude     | $\overline{A}$ | $\ast$ | 1.0     |
| $\langle r:$ delay>       | Delay         | $t_{\rm d}$    | S      | 0.0     |

<span id="page-28-4"></span>Table 12: Source type common parameters. <sup>∗</sup> depends on type.

# <span id="page-28-2"></span>4.3 Distributed "soft" current source types

```
EX <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: type> \
   <t: wfName> [ <r: size> [ <r: delay> ] ]
```


<span id="page-29-1"></span>Figure 8: Spatial region occupied by a soft electric current density source.

"Soft sources" are added to the fields already on the computational grid and as such do not cause reflections of waves scattered back towards them. Formally all such sources are ideal current sources with infinite impedance, presenting an open-circuit to any waves incident upon them. The field at the source location is the superposition of the incident field from the impressed current and the scattered field (which may be the reaction of the grid to the incident field of the source itself or from another source in the grid). The distributed soft sources supported are shown in Table [13.](#page-30-1) For surface current densities the first coordinate direction in the tag denotes the polarisation direction of the current and the second denotes the other tangential direction of the current sheet. So JMSYZ would denote a magnetic surface current density on an x-normal surface polarised in the y-direction. For all other types the single coordinate direction in the tag denotes the polarisation, so JY would denote a y-polarised electric current density.

Electric current sources are excited on primary grid edges and magnetic current sources on secondary grid edges. Figure [8](#page-29-1) shows the case of a z-polarised electric current density source on edge  $[i, i, j, j, k, k + 1]$ . Note that the electric current is evaluated at half-intergral times  $(n+\frac{1}{2})\Delta t$ ; magnetic currents are evaluated at integral times n∆t. For sources distributed over more than one edge the current density is distributed evenly over the cross-sectional area of the source using parallel current sources and uniformly along the polarisation direction of the source using equal series current sources. Essentially all the soft current sources excite the grid in the same way - the only difference is the normalisation, and units, of the amplitude of the source. Depending on the type of source been modelled one choice may be more natural than the other. If all outputs are referenced to a calibration value measured on the computational grid then the amplitude normalisation is unimportant. The normalisations for z-polarised electric current sources are

$$
J_z^{n+\frac{1}{2}}(i,j,k+\frac{1}{2}) = \mathbf{JZ} = \frac{\mathbf{JSZX}}{\tilde{\Delta}y_j} = \frac{\mathbf{JSZY}}{\tilde{\Delta}x_i} = \frac{\mathbf{IZ}}{\tilde{\Delta}x_i\tilde{\Delta}y_j} = \frac{\mathbf{I}\mathbf{DZ}}{\tilde{\Delta}x_i\tilde{\Delta}y_j\Delta z_k} = -\frac{\mathbf{EZ}}{\eta_0c_0\Delta t} = \frac{\mathbf{VZ}}{\eta_0c_0\Delta t\Delta z_k}.
$$
 (4.11)

For an x-polarised magnetic current source the equivalent normalisation is

$$
J_{\text{Mx}}^{n+\frac{1}{2}}(i,j+\frac{1}{2},k+\frac{1}{2}) = \text{JMX} = \frac{\text{JMSXY}}{\Delta z_k} = \frac{\text{JMSXZ}}{\Delta y_j} = \frac{\text{IMX}}{\Delta y_j \Delta z_k} = \frac{\text{IMDX}}{\tilde{\Delta} x_i \Delta y_j \Delta z_k} = -\frac{\eta_0 \text{HX}}{c_0 \Delta t}.
$$
(4.12)

In the above, the normalisations for the fields and voltage assume that the source is imposed in free-space.

Note that isolated current sources in free-space often lead to permanent DC charging of the computational grid due to the lack of charge dynamics in the FDTD method. As a waveform is impressed by a current source the divergent fields implicitly displace charge between cells in the grid according to the Gauss' Laws for electric and magnetic fields. This charging will be temporary if the waveform has no DC component, however, for waveforms with DC components (like a Gaussian pulse) a permanent static field will be left on the comptutational grid. If there is a conductive path between the end-points of the current source then the charge can discharge with a time constant determined by the grid capacitance and conductance.

#### <span id="page-29-0"></span>4.3.1 JX, JY, JZ, JMX, JMY and JMZ volume current density sources

Volume current densities are applied directly to grid edges - primary grid edges for electric current densities and secondary grid edges for magnetic current densities. They can be regarded as existing across an area

| Source type                              | Description                       | Units                         |
|--|-----------------------------------|-------------------------------|
| JX, JY, JZ                               | Electric (volume) current density | $\rm A\,m^{-2}$               |
| JMX, JMY, JMZ                            | Magnetic (volume) current density | $\mathrm{V}\,\mathrm{m}^{-2}$ |
| JSXY, JSXZ, JSYZ, JSYX, JSZX, JSZY       | Electric surface current density  | $Am^{-1}$                     |
| JMSXY, JMSXZ, JMSYZ, JMSYX, JMSZX, JMSZY | Magnetic surface current density  | $\rm V\,m^{-1}$               |
| IX, IY, IZ                               | Electric (line) current           | $\mathbf{A}$                  |
| IMX, IMY, IMZ                            | Magnetic (line) current           | V                             |
| IDX, IDY, IDZ                            | Electric current moment           | Am                            |
| IMDX, IMDY, IMDZ                         | Magnetic current moment           | V <sub>m</sub>                |
| VX, VY, VZ                               | Voltage                           | $\mathbf{V}$                  |
| EX, EY, EZ                               | Electric field                    | $\mathrm{V}\,\mathrm{m}^{-1}$ |
| HX, HY, HZ                               | Magnetic field                    | $\rm A\,m^{-1}$               |

<span id="page-30-1"></span>Table 13: Distributed "soft" current source types.  $1 \text{ Wb} = 1 \text{ Vs}$ .

normal to their polarisation ditrection given by the area of the dual grid faces associated with the edges excited. For example,

# A soft electric current density source EX 16 17 16 17 16 17 current\_density JY wf1 1.0

would correspond to a y-polarised electric current density of  $1 \text{ A m}^{-2}$  impressed on four primary grid edges in the volume bounded by [15.5,17.5,15.5,17.5,16,17].

# <span id="page-30-0"></span>4.3.2 JSXY, JSXZ, JSYX, JSYZ, JSZX, JSZY, JMSXY, JMSXZ, JMSYX, JMSYZ, JMSZX, JMSZY surface current density sources

Surface current density sources are useful for impressing "current sheet" sources across surfaces of the grid. According to Love's Surface Equivalence Principle the sources within a surface S can be replaced by the equivalent surface currents

$$
\mathbf{J}_{\mathrm{Ms}} = -\hat{\mathbf{n}} \times \mathbf{E}|_{S} \tag{4.13}
$$

$$
\mathbf{J}_{\mathrm{s}} = \hat{\mathbf{n}} \times \mathbf{H}|_{S} \tag{4.14}
$$

on the surface, leaving a null field inside the surface and the original field outside the surface. An aperture in a PEC surface for example can be modelled using a magnetic surface current density  $J_{\text{Ms}} = -\hat{\boldsymbol{n}} \times \boldsymbol{E}|_{\text{s}}$ radiating on a PEC surface with the aperture closed (the electric current density is shorted out in this case). The known electric field in an aperture can be impressed using

```
# An equivalent magnetic current density source for an
# aperture with an electric field of 1 V/m.
TB 0 10 10 10 0 10 PEC
EX 5 6 10 10 5 6 aperture_efield JMSZY wf1 1.0
```
Note that this will radiate into both the  $j < 10$  and  $j > 10$  half-spaces which is probably not what is intended. A volumetic PEC backing needs to be used on the non-radiating side, say the  $j < 10$  half-space:

```
# An equivalent magnetic current density source for an
# aperture with a uniform electric field of 1 V/m.
MB 0 10 0 10 0 10 PEC
EX 5 6 10 10 5 6 aperture_efield JMSZY wf1 1.0
```
Another example is

# A soft electric surface current density source EX 15 20 15 20 10 10 current\_sheet JSXY wf1 1.0

which corresponds to an x-polarised electric surface current density of  $1 \text{ A m}^{-1}$  impressed on 36 primary grid edges on the z-normal surface bounded by [15,20,14.5,20.5,9.5,10.5]. The equivalent volume current density is determined by dividing surface current density by the dual grid edge length normal to the surface, in this case  $\tilde{\Delta}z_{10}$ .

For infinitesimal sources that are not aligned with the grid it may be better to distributed the source over two edges for each polarisation so that each polarisation has the same phase centre [\[6\]](#page-66-3). For example and over two eages for each polarisation so that each polarisation has the same phase centre  $[0]$ . For example an<br>electric dipole moment with amplitude 1 A m, aligned along unit vector  $(1\sqrt{3}, 1\sqrt{3}, 1\sqrt{3})$  can be implemen using

```
# Non-aligned source with same phase centre for each polarisation.
EX 10 11 10 10 10 10 current_sheet IDX wf1 0.5773
EX 10 10 10 11 10 10 current_sheet IDY wf1 0.5773
EX 10 10 10 10 10 11 current_sheet IDZ wf1 0.5773
```
# <span id="page-31-0"></span>4.3.3 IX, IY, IZ, IMX, IMY and IMZ line current sources

Line current sources are related to current densities using the area of the dual grid face normal to and centred on the excited edge. For example,

```
# A soft electric current source
EX 15 15 15 15 14 16 line_current IZ wf1 1.0
```
impresses a z-directed current of 1 A flowing in the volume [14.5,15.5,14.5,15.5,14,16]. The current density is determined as  $1/(\tilde{\Delta}x_{15}\tilde{\Delta}y_{15})\,\text{A}\,\text{m}^{-2}$ . Current sources can be specified on volumetric bounding boxes as well. So,

# A soft volumetric electric current source EX 15 16 15 16 14 16 line\_current IZ wf1 1.0

would distribute the 1 A current across the area  $(\tilde{\Delta}x_{15} + \tilde{\Delta}x_{16}) \times (\tilde{\Delta}y_{15} + \tilde{\Delta}y_{16})$  normal to the current direction.

# <span id="page-31-1"></span>4.3.4 IDX, IDY, IDZ, IMDX, IMDY and IMDZ current moment sources

Current moments impress the specified current moment on a volume, surface or edge. The current is determined by dividing the moment by the total length of the bounding box in the polarisation direction and then the current density is found by dividing by the area of the normal dual grid faces. For example,

# # A soft volumetric electric current moment source EX 15 16 15 16 14 16 current\_moment IDZ wf1 1.0

would distribute the 1 A m current moment across the area  $(\tilde{\Delta}x_{15} + \tilde{\Delta}x_{16}) \times (\tilde{\Delta}y_{15} + \tilde{\Delta}y_{16})$  normal to the current direction and length  $\Delta z_{14} + \Delta x_{15}$ .

# <span id="page-31-2"></span>4.3.5 EX, EY, EZ, HX, HY and HZ soft "field" sources

Soft "field" sources are added directly to the fields already on the grid, however, as noted above they are formally current sources with respect to their physical behaviour. The field they impress on the grid is dependent on the time-step (and hence CFLN). In general this means they are difficult to use on a nonuniform mesh. In Vulture the soft electric field source is normalised such that on a uniform grid with the uniform mesn. In vulture the soft electric field source is normalised such that on a uniform grid with the default CFLN of  $\sqrt{3}/2$  an infinite plane of colinearly polarised edges would launch a plane wave with the specified electric/magnetic field amplitude in to the grid. Explicitly,

```
# A soft electric ''field'' source
EX 1 10 1 10 5 5 eplane EZ wf1 1.0
```
would launch a "plane-wave" of amplitude 1 V/m in both the positve and negative z-directions on a uniform mesh with the default CFLN. Note that in practice the source plane is usually not infinite and therefore this give a poor representation of a plane-wave: There will be always be diffraction from the edges of the source plane. However, within an infinite parallel-plate waveguide, simulated using PEC and PMC external boundaries, the source plane is effectively infinite and in this case a well defined TEM wave is launched (in both directions). The PEC external boundaries also effectively remove any displaced charge from the grid and prevent DC charging of the grid for this case too.

# <span id="page-31-3"></span>4.3.6 VX, VY and VZ soft "voltage" sources

Soft voltage sources are just soft field sources normalised by the total edge length of the bounding box in the direction of polarisation, so all the comments above for field sources apply equally here.

# <span id="page-32-0"></span>4.4 Distributed "hard" field source types

EX <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: type> \ <t: wfName> [ <r: size> [ <r: delay> ] ]



<span id="page-32-1"></span>Figure 9: Spatial region occupied by a hard electric field source.

"Hard sources" replace the fields already on the computational grid and as such cause reflections of waves scattered back towards them. Formally such sources are ideal *voltage sources* with zero impedance, presenting a short-circuit to any waves incident upon them. For sources distributed over more than one edge the voltage is distributed uniformly over the cross-sectional area of the source using equal parallel voltage sources and evenly along the polarisation direction of the source using series voltage sources. The edges on which hard sources are impressed appear as PEC or PMC materials as far as waves scattering from them are concerned. Indeed, one practical use of a hard source is to impress the equivalent surface currents of an aperture field onto a PEC surface, however, it must always be bourne in mind that any waves reflected back towards such a source may suffer unphysical reflections. This is discussed futher below. The distributed hard sources supported are shown in Table [13.](#page-30-1)

Electric sources are excited on primary grid edges (centres of secondary grid faces) and magnetic sources at the centres of primary grid faces (on secondary grid edges). Figure [9](#page-32-1) shows the case of a z-polarised hard electric field source on edge  $[i, i, j, j, k, k + 1]$ . Note that electric field sources are evaluated at intergral times  $n\Delta t$ ; magnetic field sources are evaluated at half-integral times  $(n + \frac{1}{2})\Delta t$ . Essentially all the hard field sources excite the grid in the same way - the only difference is the normalisation, and units, of the amplitude of the source. Depending on the type of source been modelled one choice may be more natural than the other. If all outputs are referenced to a calibration value measure on the computational grid then the amplitude normalisation is unimportant. The normalisations for a z-polarised hard electric field source are

$$
E_z^{n+\frac{1}{2}}(i,j,k+\frac{1}{2}) = \mathbf{EZ} = -\frac{\mathbf{VZ}}{\Delta z_k} = -\eta_0 c_0 \Delta t \mathbf{JZ} = -\eta_0 c_0 \Delta t \frac{\mathbf{JSZX}}{\tilde{\Delta}y_j} = -\eta_0 c_0 \Delta t \frac{\mathbf{JSZX}}{\tilde{\Delta}x_i}
$$
  

$$
= -\eta_0 c_0 \Delta t \frac{\mathbf{IZ}}{\tilde{\Delta}x_i \tilde{\Delta}y_j} = -\eta_0 c_0 \Delta t \frac{\mathbf{IZZ}}{\tilde{\Delta}x_i \tilde{\Delta}y_j \Delta z_k}
$$
(4.15)

For an x-polarisaed hard magnetic field source the equivalent normalisation is

$$
H_{\mathbf{x}}^{n+\frac{1}{2}}(i, j+\frac{1}{2}, k+\frac{1}{2}) = \mathbf{H}\mathbf{x} = -\frac{c_0 \Delta t}{\eta_0} \mathbf{J} \mathbf{M} \mathbf{x} = -\frac{c_0 \Delta t}{\eta_0} \frac{\mathbf{J} \mathbf{M} \mathbf{S} \mathbf{X} \mathbf{Y}}{\Delta z_k} = -\frac{c_0 \Delta t}{\eta_0} \frac{\mathbf{J} \mathbf{M} \mathbf{S} \mathbf{X} \mathbf{Z}}{\Delta y_j}
$$

$$
= -\frac{c_0 \Delta t}{\eta_0} \frac{\mathbf{I} \mathbf{M} \mathbf{X}}{\Delta y_j \Delta z_k} = -\frac{c_0 \Delta t}{\eta_0} \frac{\mathbf{I} \mathbf{M} \mathbf{D} \mathbf{X}}{\tilde{\Delta} x_i \Delta y_j \Delta z_k}.
$$
(4.16)

The above the normalisations for the fields and voltage assume that the source is imposed in free-space. Note that the fields induced on the grid for all the current related hard sources depend on the time-step (and CFLN).

| Source type   | Description                       | Units           |
|---|-----------------------------------|-----------------|
| $=EX$ , $=EY$ , $=EZ$   | Electric field                    | $\rm V\,m^{-1}$ |
| $=$ HX, $=$ HY, $=$ HZ  | Magnetic field                    | $Am^{-1}$       |
| $=VX$ , $=VY$ , $=VZ$   | Voltage                           | V               |
| $=$ JX, $=$ JY, $=$ JZ  | Electric (volume) current density | $Am^{-2}$       |
| $=JMX$ , $=JMY$ , $=JMZ$  | Magnetic (volume) current density | $\rm V\,m^{-2}$ |
| $=$ JSXY, $=$ JSXZ, $=$ JSYZ, $=$ JSYX, $=$ JSZX, $=$ JSZY      | Electric surface current density  | $Am^{-1}$       |
| $=JMSXY$ , $=JMSXZ$ , $=JMSYZ$ , $=JMSYX$ , $=JMSZX$ , $=JMSZY$ | Magnetic surface current density  | $\rm V\,m^{-1}$ |
| $=IX, =IY, =IZ$   | Electric (line) current           | $\mathbf{A}$    |
| $=IMX$ , $=IMY$ , $=IMZ$  | Magnetic (line) current           | V               |
| $=IDX$ , $=IDY$ , $=IDZ$  | Electric current moment           | A m             |
| $=IMDX$ , $=IMDY$ , $=IMDZ$                                     | Magnetic current moment           | V <sub>m</sub>  |

Table 14: Distributed "hard" field source types.

#### <span id="page-33-0"></span>4.4.1 =  $EX$ ,  $= EY$ ,  $= EZ$ ,  $= HX$ ,  $= HY$  and  $= HZ$  hard field sources

Hard field source can be used to impress known field distributions on surfaces, similar to the use of current sheets in Section [4.3.2](#page-30-0) above. With a hard source an aperture field can be imposed directly on a PEC surface, for example:

```
# Direct Dirchlet boundary condition for
# aperture with a uniform electric field of 1 V/m.
MB 0 10 0 10 0 10 PEC
EX 5 6 10 10 5 6 aperture_efield =EZ wf1 1.0
```
#### <span id="page-33-1"></span>4.4.2 =VX, =VY and =VZ hard voltage sources

Hard voltage sources are just hard field sources normalised by the total edge length of the bounding box in the direction of polarisation. They can be used in cases where an ideal voltage source excitation is required.

### <span id="page-33-2"></span>4.4.3 =JX, =JY, =JZ, =JMX, =JMY and =JMZ hard current density sources

These source types are provided for completeness but may not be practically useful.

### <span id="page-33-3"></span>4.4.4 =JSXY, =JSXZ, =JSYX, =JSYZ, =JSZX, =JSZY, =JMSXY, =JMSXZ, =JMSYX, =JMSYZ, =JMSZX, =JMSZY hard surface current density sources

These source types can be used to impose current sheets, however, the field induced on the mesh is dependent on the time-step and will only have the correct normalisation on uniform cubic grids with the default CFLN. It is usually better to imposed field values directly using =EX etc.

#### <span id="page-33-4"></span>4.4.5 =IX, =IY, =IZ, =IMX, =IMY and =IMZ hard current sources

These source types are provided for completeness but may not be practically useful.

#### <span id="page-33-5"></span>4.4.6 =IDX, =IDY, =IDZ, =IMDX, =IMDY and =IMDZ hard current moment sources

These source types are provided for completeness but may not be practically useful.

# <span id="page-33-6"></span>4.5 Other source types

#### <span id="page-33-7"></span>4.5.1 VRX, VRY, VRZ, IRX, IRY and IRZ finite impedance lumped sources

EX <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> VRX|VRY|VRZ|IRX|IRY|IRZ \ <t: wfName> <r: resist> [ <r: size> [ <r: delay> ] ]

This form of the EX directive is used to introduce lumped Thévenin voltage and Norton current sources with finite internal impedance. For Thévenin voltage sources the amplitude  $\langle r: \text{ size} \rangle$  is the open circuit voltage of the source in volts and  $\leq r$ : resist> is the internal resistance in ohms. For Norton current sources



<span id="page-34-2"></span>Figure 10: Schematic illustration of lumped Thévenin voltage source (left) and Norton current source (right) on a z-polarised edge.

the amplitude  $\langle r: \text{ size} \rangle$  is the short circuit current of the source in amps and  $\langle r: \text{ resist} \rangle$  is the internal resistance in ohms. Both sources are implement as soft current sources on the corresponding electric field edge, together with an appropriate conductive loading of the edge to account for the source resistance, as shown in Figure [10.](#page-34-2) Currently the bounding box must be a single edge!

Examples:

```
# A soft z-polarised lumped voltage source of ampltiude 1 V
# and internal resistance 50 ohms applied on the grid edge
# from (16,16,16) to (16,16,17).
EX 16 16 16 16 16 17 feed VRZ wf1 50.0 1.0
```
<span id="page-34-0"></span>4.5.2 VDX, VDY and VDZ delta-gap thin wire voltage sources

Delta-gap sources are currently not implemented!

# <span id="page-34-1"></span>4.6 Plane-wave sources: PW

Plane-waves can be launched into the computational grid using the PW directive:

```
PW <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: wfName> \
   <r: theta> <r: phi> <r: psi> [ <s: mask> [<r: size> [ <r: delay> ] ] ]
```
Plane-wave sources partition the grid into two regions:

- A total field region inside of the bounding box.
- A scattered field region outside of the bounding box.

Such source are also known as Huygen's sources or total-field scattered-field (TFSF) sources or boundaries. The bounding box must define a volume and *must not cut across any objects on the mesh*; it can however butt-up against an internal PEC surface or external PEC, PMC or PERIODIC surface in certain cases (see <t: mask> below). Plane-wave sources are applied in the order they are given in the mesh file after all other types of sources.

| Parameter                        | Quantity           | Symbol      | Units   | Default |
|----------------------------------|--------------------|-------------|---------|---------|
| lt: name                         | Plane-wave name    | n/a         | n/a     | none    |
| $lt:$ wfName>                    | Waveform name      | n/a         | n/a     | none    |
| $\langle r: \text{theta}\rangle$ | Polar angle        | $\theta$    | degrees | none    |
| $\langle r: \text{phi}\rangle$   | Azimuthal angle    | $\phi$      | degrees | none    |
| $\langle r: \text{psi}\rangle$   | Polarisation angle | ψ           | degrees | none    |
| $\ltt$ : mask>                   | Face activity mask | n/a         | n/a     | 111111  |
| $\langle r: size \rangle$        | Magnitude          | $E_0$       | V/m     | 1.0     |
| $\langle r:$ delay>              | Delay              | $t_{\rm d}$ | S       | 0.0     |

<span id="page-34-3"></span>Table 15: Plane-wave source type parameters.

The plane-wave source type parameters are:

- <t: name>: The unique name of the plane-wave.
- $\leq$ t: wfName>: The name of the waveform to use for the source. This must have been defined already in a WF directive.
- $\leq r$ : theta>: The polar angle,  $0 \leq \theta \leq 180$ , of the wave vector in degrees.
- $\leq r$ : phi>: The azimuthal angle,  $0 \leq \phi < 360$ , of the wave vector in degrees.
- $\langle \mathbf{r}: \mathsf{psi} \rangle$ : The polarisation angle,  $0 \leq \psi \leq 360$ , of the electric field vector relative to  $\hat{\mathbf{k}}^{\text{inc}} \times \hat{\mathbf{z}}$  in degrees.
- <s: mask>: A bit mask of six binary digits indicating which faces (XLO XHI YLO YHI ZLO ZHI) of the plane-wave bounding box are active. e.g 111111 indicates all faces are active while 011111 indicates that the XLO face is inactive. This can be used to terminate one face (the inactive one) of the TFSF surface with a PEC plate to construct a semi-infinite space. The reflected wave must be included using a second PW directive with opposite wave-vector and polarisation in this case to obtain accurate results - see the examples later.
- $\langle \mathbf{r}: \text{ size}\rangle$ : The source amplitude,  $E_0$ , in V/m. The default value is 1.0 V/m.
- $\leq$  r: delay>: The delay of the source relative to the waveform in seconds. This delay is in addition to any delay specified in the WF directive.

These parameters are summarised in Table [15](#page-34-3) and Figure [11](#page-35-0) shows how the angles are defined. The incident wave-vector, electric field vector and magnetic field vector are given by

$$
\hat{\mathbf{k}}^{\text{inc}} = \hat{\mathbf{r}} = \begin{bmatrix} \sin(\theta)\cos(\phi) \\ \sin(\theta)\sin(\phi) \\ \cos(\theta) \end{bmatrix}
$$
\n(4.17)

$$
\mathbf{E}^{\rm inc} = E_0 \begin{bmatrix} \cos(\psi)\sin(\phi) - \sin(\psi)\cos(\theta)\cos(\phi) \\ -\cos(\psi)\cos(\phi) - \sin(\psi)\cos(\theta)\sin(\phi) \\ \sin(\psi)\sin(\theta) \end{bmatrix}
$$
(4.18)

$$
\mathbf{H}^{\rm inc} = \frac{E_0}{\eta_0} \begin{bmatrix} \sin(\psi)\sin(\phi) + \cos(\psi)\cos(\theta)\cos(\phi) \\ -\sin(\psi)\cos(\phi) + \cos(\psi)\cos(\theta)\sin(\phi) \\ -\cos(\psi)\sin(\theta) \end{bmatrix} . \tag{4.19}
$$

The angles for common directions and polarisations are given in Table [16.](#page-36-4)



<span id="page-35-0"></span>Figure 11: Definition of the angles of the wave-vector and electric field vector for the plane-wave source.

On a cubic mesh an auxiliary 1D grid is used to accurately account for dispersion across the total field region. On non-cubic uniform meshes and non-uniform meshes a basic analytic incident field is used: This has low accuracy on electrically large meshes. In any case it is advisable to make the plane-wave bounding box as small as possible to minimise any dispersion errors.
| $\hat{\mathbf{k}}^{\mathrm{inc}}$ | $\hat{\mathbf{E}}^{\text{inc}}$ | $\mathbf{H}^{\text{inc}}$ | $\theta$ (° | $\phi(^{\circ})$ | $\psi$ (°) |
|-----------------------------------|---------------------------------|---------------------------|-------------|------------------|------------|
| $k_x\hat{\textbf{x}}$             | $E_y \hat{\textbf{y}}$          | $H_z\hat{\mathbf{z}}$     | 90          | 0                | 180        |
| $k_x\hat{\textbf{x}}$             | $E_z\hat{\mathbf{z}}$           | $H_y \hat{\textbf{y}}$    | 90          | 0                | 90         |
| $k_y \hat{\mathbf{y}}$            | $E_x\hat{\textbf{x}}$           | $H_z\hat{\mathbf{z}}$     | 90          | 90               | 0          |
| $k_y \hat{\mathbf{y}}$            | $E_z\hat{\mathbf{z}}$           | $H_x\hat{\textbf{x}}$     | 90          | 90               | 90         |
| $k_z \hat{\mathbf{z}}$            | $E_x\hat{\textbf{x}}$           | $H_y \hat{\textbf{y}}$    | 0           | 0                | 270        |
| $k_z \hat{\mathbf{z}}$            | $E_y \hat{\textbf{y}}$          | $H_x\hat{\textbf{x}}$     | 0           | 0                | 180        |
| $-k_x\hat{\mathbf{x}}$            | $-E_y\hat{\textbf{y}}$          | $H_z\hat{\mathbf{z}}$     | 90          | 180              | 180        |
| $-k_x\hat{\mathbf{x}}$            | $-E_z\hat{\mathbf{z}}$          | $-H_y\hat{\textbf{y}}$    | 90          | 180              | 270        |
| $-k_y\hat{\textbf{y}}$            | $-E_x\hat{\mathbf{x}}$          | $-H_z\hat{\mathbf{z}}$    | 90          | 270              | 0          |
| $-k_y\hat{\textbf{y}}$            | $-E_z\hat{\mathbf{z}}$          | $H_x\hat{\textbf{x}}$     | 90          | 270              | 270        |
| $-k_z\hat{z}$                     | $-E_x\hat{\mathbf{x}}$          | $H_y \hat{\textbf{y}}$    | 180         | 0                | 270        |
| $-k_z\hat{\mathbf{z}}$            | $-E_y\hat{\textbf{y}}$          | $-H_x\hat{\mathbf{x}}$    | 180         | 0                | 0          |

Table 16: Angles for common directions of propagation and polarisation.

# 5 Observers

The observer directive format and assocaited output file formats are under review and may change in a future version of Vulture.

# 5.1 Simulation output control: OT, OF

The OT directive determines for which time-steps data are output by the time-domain observers:

OT <i: tstart> <i: tstop>

The data for all time-steps between and including  $\langle i: \text{start}\rangle$  and  $\langle i: \text{start}\rangle$  will be written out. By default the data for all time-steps is output if no OT directive is given.

For frequency domain observers, including waveforms and far-fields calculations, the frequencies to be output are determined by the OF directive:

OF <r: fstart> <r: fstop> <i: numFreq>

By default, if no OF directive is given, the frequencies

$$
f_k = \frac{k}{N_s \Delta t} \quad k = 0, \dots, \frac{N_s}{10}
$$
\n
$$
(5.1)
$$

are output, where  $N_s$  is the number of time-steps and  $\Delta t$  is the time-step. The DFTs are computed using the time-series within the bounds specified by any OT directive or the full time series if there is no OT directive.

## 5.2 Implicit waveform observers

For every waveform directive found in the mesh file two implicit "'waveform observers" are automatically created which monitor the waveform's time-series and frequency spectrum. Waveform time-series are output into ASCII files called  $wf\_t$ : name> $_t$ d.asc, where  $\lt t$ : name> is the name of the waveform in the meshfile. The associated frequency spectra are similarly stored in ASCII files called  $wf_{\text{-}}\text{ }ct: \text{ name}_{\text{-}}\text{ }fd$ . asc. The output format are given in Tabe [17.](#page-37-0)

# 5.3 Near-field observers: OP

Near-field output requests are made using OP directives which have two forms depending on the type:

```
OP <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: type> \
   [ \ltt: wfName> ]
OP <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> <t: type> \
   [ <i: xstep> <i: ystep> <i: zstep> ] ]
```
These directives place observers on the computational grid that output the fields on and within the specified bounding box.

```
# Waveform# 0
# ts (-) t (s) wf (-)<br>0 0.00000000e+00 0.0000000e+00
          0.00000000e+00. .............. ..............
# Waveform# 0
\#\qquad f (Hz) Re(yf) (-) Im(yf) (-)
 0.00000000e+00 0.00000000e+00 0.00000000e+00
  .............. .............. ..............
```
<span id="page-37-0"></span>Table 17: Output formats for implicit waveform obserbvers: Time series (top) and frequency spectra (bottom).

#### 5.3.1 TDOM BINARY type observers

```
OP <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t:name> TDOM_BINARY \
   [ <i: xstep> <i: ystep> <i: zstep> ]
```
The TDOM BINARY type requests time-domain field outputs over a volumetric region in the University of York finite methods binary output format [\[3\]](#page-66-0). The electric and magnetic field components in each cell enclosed by the bounding box are output with a cell index stride length in each coordinate direction given by <i: xstep>, <i: ystep> and <i: zstep>. The default stride lengths are one in all directions. The time-series of the first waveform defined in the mesh file will be stored in the excite.dat to be used as a reference in frequency domain calculations.

The observer outputs are indexed by cell numbers but are sampled at their respective positions and times in the Yee cell shown in Figure [2.](#page-12-0) For example  $\text{Ex}[i][j][k]$  is sampled at position  $(x_i + \frac{dx_i}{2}, y_j, z_k)$ and  $\text{Hx[i][j][k]}$  is sampled at position  $(x_i, y_j + \frac{dy_j}{2}, z_k + \frac{dz_k}{2})$ . Time domain fields are output with electric field times. The magnetic fields are half a time step later than the time specified.

Frequency domain outputs are divided by the spectrum of the (complex) reference waveform, defined by the <t: wfName> parameter, to give the frequency response of the observerable relative to the source waveform. Any amplitude factor in the waveform directive is therefore divided out, however, amplitude factors in excitation directives are retained. The default reference waveform is the first one found in mesh. Magnetic field spectra have a phase shift (advance) corresponding to half a time step.

The results are stored in a binary output file called impulse.dat. Results from multiple TDOM BINARY types observers are stored in the same impulse.dat file. Two other files are also created when TDOM BINARY type observers are present:

- excite.dat: An ASCII file containing the reference excitation waveform time series.
- process.dat: A template ASCII file for controlling other post processing tools.

See Section [7.2](#page-41-0) for information on using the binary output data. Note that the name tag is currently not propagated in to the output data files.

For example the directive

```
OP 4 10 14 20 10 10 output1 TDOM_BINARY 2 2 1
```
will output electric and magnetic fields on a 2D grid of cells with

- $x$  indices 4, 6, 8 and 10.
- y indices 14, 16, 18 and 20.
- $z$  index 10.

#### 5.3.2 TDOM ASCII and FDOM ASCII type observers

```
OP <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> TDOM_ASCII
OP <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> FDOM_ASCII [ <t: wfName> ]
```
The TDOM ASCII and FDOM ASCII observer types requests time-domain and frequency-domain ASCII format field outputs respectively in a single mesh cell. As such the bounding box must be a single cell.

Time-domain outputs for each observer of type TDOM ASCII are written to separate files with names eh <t: name> td.asc. The file format is given in Table [18.](#page-38-0) The outputs are sampled at their respective positions and times in the Yee cell shown in Figure [2.](#page-12-0) For example Ex[i][j][k] is sampled at position  $(x_i + \frac{dx_i}{2}, y_j, z_k)$  and Hx[i][j][k] is sampled at position  $(x_i, y_j + \frac{dy_j}{2}, z_k + \frac{dz_k}{2})$ . Time domain fields are output with electric field times. The magnetic fields are half a time step later than the time specified.

Frequency-domain outputs for each observer of type FDOM ASCII are written to separate files with names eh\_<t: name>\_fd,asc. The file format is given in Table [19.](#page-38-1) Frequency domain outputs are divided by the spectrum of the (complex) reference waveform, defined by the  $\leq t$ : wfName> parameter, to give the frequency response of the observerable relative to the source waveform. Any amplitude factor in the waveform directive is therefore divided out, however, amplitude factors in excitation directives are retained. If no reference waveform is given the first one found in the mesh file is used. Magnetic field spectra have a phase shift corresponding to half a time step.

```
# [<i: ilo>,<i: jlo>,<i: klo>] -> (x_ilo,y_jlo,z_klo)
# ts (-) t (s) Ex (V/m) Ey (V/m) Ez (V/m) Hx (A/m) ...
     0 0.00000000e+00 0.00000000e+00 0.00000000e+00 0.00000000e+00 0.00000000e+00 ...
         . .............. .............. .............. .............. .............. ...
```
<span id="page-38-0"></span>Table 18: Output format of the TDOM ASCII observer types.



<span id="page-38-1"></span>Table 19: Output format of the FDOM ASCII observer types.

#### 5.3.3 HDF5 type observers

HDF5 observers are currently not implemented!

OP <i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <t: name> TDOM\_HDF5 [<t: wfName>]

These will have the following characteristics:

- Nodal observers.
- All outputs to one file called output.h5.
- Maybe AMELET format.
- Optional compilation using USE\_HDF5=0N/OFF.

## 5.4 Far-field observers: FF

Far-field observers are currently not implemented!

The FF directive requests a near-field to far-field transformation of the fields over the bounding box:

FF  $\langle i: ilo \rangle \langle i: ihi \rangle \langle i: jlo \rangle \langle i: jhi \rangle \langle i: kho \rangle \langle i: khi \rangle \langle t: name \rangle \langle t: wfName \rangle$ <r: theta1> <r: theta2> <i: num\_theta> <r: phi1> <r: phi2> <i: num\_phi> [ <t: mask> ]

These will have the following characteristics:

- Calculate the frequency domain far-field from the near-field sampled on the bounding box.
- Calculate the total radiated power in far field by integrating the pattern.
- Calculate the frequency domain power flowing out of the bounding box by integrating the Poynting flux.
- Check the power balance between the previous two powers.
- Use a running DFT via "internal" surface observers.
- Use frequencies from the OF card or those automatically determined.
- The bounding box must be a volume.
- $\leq r$ : theta1>: Start polar angle in degrees [0,180).
- $\langle r: \text{theta2}\rangle$ : Stop polar angle in degrees  $(0,180), \geq r: \text{theta1}$
- $\leq i$ : num theta>: Number of polar angles,  $\geq 0$ .
- $\langle \mathbf{r}: \text{phi12}: \text{Start azimuthal angle in degrees } [0,360)$ .
- $\langle \mathbf{r}: \text{phi2}\rangle$ : Stop azimuthal angle in degrees  $(0,360], \rangle = \mathbf{r}: \text{phi1}.$
- $\leq i$ : num phi>: Number of azimuthal angles,  $\geq 0$ .
- <t: mask> is a bitmap that defines which faces (XLO XHI YLO YHI ZLO ZHI) are included e.g 111111, 011111.
- Origin is the centre of bounding box may need to make it a parameter?
- Phase reference is the waveform? Distance for phase calculation?
- Output is  $\theta$ ,  $\phi$ ,  $E_{\theta}$ ,  $E_{\phi}$ , power gain/RCS in files called rcs <i: freqNum>.asc where <i: freqNum> is the frequency number.

# 6 Viewing the mesh

## 6.1 gvulture

The program gvulture converts a Vulture mesh into a set of files suitable for 3D viewing of the mesh using Gnuplot [\[7\]](#page-66-1) or Gmsh [\[8\]](#page-66-2). The gvulture program takes a Vulture mesh file as input and writes out a number of files to the current directory.

A typical command line for running gvulture is

#### \$ gvulture -p antenna.mesh

where **antenna**.mesh is the name of the input mesh file and the  $-p$  option requests the output to be in physical mesh units rather than the default which uses the mesh line indices as coordinates. The default output format is for Gnuplot. To generate Gmsh output format add the -m flag:

#### \$ gvulture -p -m antenna.mesh

gvulture also creates a log file called gvulture.log. To suppress the inclusion of representations of the external mesh surfaces in the output files use the -e options:

#### \$ gvulture -p -m -e antenna.mesh

Further specific details on the two output format are given in the following sub-sections.

# 6.2 Gnuplot

Gnuplot is public domain graph plotting program that will run on most display devices [\[7\]](#page-66-1). It can also produce (Encapsulated) Postscript, Latex picture environment and various bitmap image format outputs. When used in the default mode to generate Gnuplot output gvulture generates a number of files:

- mesh.gnp: An ASCII Gnuplot script that must be executed in either batch or interactive mode to create the image of the mesh.
- gnuplot-lines.dat: An ASCII data file containing a Gnuplot compatible description of the mesh lines.
- gnuplot-external.dat: An ASCII data file containing a Gnuplot compatible description of the external surfaces of the mesh. The edges of the cuboid forming the simulation space are shown.
- gnuplot-surface.dat: An ASCII data file containing a Gnuplot compatible description of the internal surfaces in the mesh. The outer perimeter of the surfaces are shown.
- gnuplot-blocks.dat: An ASCII data file containing a Gnuplot compatible description of the materials blocks in the mesh. The outer surface of the volume of the material block are shown.
- gnuplot-sources.dat: An ASCII data file containing a Gnuplot compatible description of the sources in the mesh. These are shown as the bounding box of the sources with an arrow indicating the direction of the electric field polarisation.
- gnuplot-planewaves.dat: An ASCII data file containing a Gnuplot compatible description of the plane-waves in the mesh. These are shown as the bounding boxes of the active faces of the plane-wave excitation surface. An orthogonal set of arrows are added to the corner of the plane-wave surface where the wavefront first enters the mesh. The direction of propagation is denoted by an open arrow, the electric field direction by a closed arrow and the magnetic field direction by a double closed arrow.
- gnuplot-observers.dat: An ASCII data file containing a Gnuplot compatible description of the observers in the mesh. These are rendered as points at the nodes where outputs are generated.

An image of the mesh from the default viewing direction can created in batch mode from the command line using

#### \$ gnuplot mesh.gnp

resulting in an Encapsulated Postscript file mesh.eps. To view the mesh interactively start gnuplot in interactive mode and use the load command to load the mesh.gnp file:

```
$ gnuplot
gnuplot> load 'mesh.gnp'
```
The viewing direction can then be changed using the mouse or by issuing commands to the Gnuplot interpreter [\[9\]](#page-66-3).

# 6.3 Gmsh

When used with the  $-m$  option to generate Gmsh format output gvulture generates a single output Gmsh ASCII format file called mesh.msh. This can be loaded into Gmsh using

#### \$ gmsh mesh.msh

Details on using the gmsh program can be found in its manual [\[10\]](#page-66-4). The most useful actions for viewing Vulture meshes include:

- Rotating and scaling the view using the mouse.
- Selecting the physical groups in the mesh that are visible.
- Altering the appearance of the mesh items that are rendered.
- Saving the current view as an image file.

# 7 Post-processing tools

# 7.1 TDOM ASCII and FDOM ASCII type outputs

These observer type have outputs that are flat ASCII files which can easily be plotted using Gnuplot or GNU Octave or imported into spreadsheets etc.

# <span id="page-41-0"></span>7.2 TDOM BINARY type outputs

## 7.2.1 Overview

A suite of programs is available for post-processing the time response stored in the impulse.dat file created by TDOM BINARY type observers. The processing can be done in either the time-domain or, via a Fast Fourier Transform (FFT), in the frequency-domain. Full details can be found in the Processing Tools Manual [\[3\]](#page-66-0).

These programs require the files impulse.dat, process.dat and excite.dat which are generated by Vulture for TDOM\_BINARY types observers. The files impulse.dat and excite.dat should be not be modified; only the file process.dat require alterations to control the output programs described below. The format of process.dat is shown in Table [20.](#page-41-1)

```
<i: nobs>
<i: ilo> <i:ihi> <i:istep> <i:jlo> <i:jhi> <i:jstep> <i:klo> <i:khi> <i:kstep>
.................
<i: ilo> <i:ihi> <i:istep> <i:jlo> <i:jhi> <i:jstep> <i:klo> <i:khi> <i:kstep>
<i: n1> <i: n2> <i: nstep>
\langle r: f1 \rangle \langle r: f2 \rangle \langle r: fstep \rangle<r: meshsize>
<i: ilo> <i: ihi> <i: jlo> <i: jhi> <i: klo> <i: khi> <i: field>
```
<span id="page-41-1"></span>Table 20: Format of the process.dat file.

The parameters in the file have the following interpretations:

- <i: nobs>: is the number of TDOM BINARY type observers. Do not modify.
- The following  $\langle i: \text{nobs} \rangle$  lines are the bounding boxes  $[\langle i: \text{ilo} \rangle, \langle i: \text{ jlo} \rangle, \langle i: \text{ jhi} \rangle]$ <i: klo>,<i: khi>] and stride lengths [<i: istep>,<i: ijstep>,<i: kstep>] of these TDOM BINARY observers in the order given in the mesh file. Do not modify.
- $\leq i: n1$ ,  $\leq i: n2$  and  $\leq i: n$  step give the start time-step, stop time-step and number of time timesteps respectively, initially as defined in the OT and NT directives. These can be modified to the time or frequency iterations for which data should be output.
- $\langle$ r: f1>,  $\langle$ r: f2> and  $\langle$ f: fstep> are the first frequency, last frequency and frequency step, initially as defined in the OF directive (or determined automatically by the solver). Do not modify.
- $\leq$  r: meshsize> is the mesh size. For non-cubic meshes this value is not meaningful the processing tools do not fully support non-uniform meshes and only work with mesh line indices. Do not modify.
- The last can be modified to give the bounding box  $\lceil \langle i : i \bigr\rceil$ ,  $\langle i : i \bigr\rceil$ ,  $\langle i : j \bigr\rceil$ ,  $\langle i : j \bigr\rceil$ ,  $\langle i : k \bigr\rceil$ ,  $\langle i : k \bigr\rceil$ ,  $\langle i : k \bigr\rceil$ and the field component  $\leq i:$  field> of the outputs to be extracted as described in the following. The meaning of the  $\leq i$ : field> parameter is defnied in Table [21.](#page-41-2)

| $\langle i: \text{field} \rangle$ | Field component |
|-----------------------------------|-----------------|
|                                   | $E_x$           |
| 2                                 | $E_y$<br>$E_z$  |
| 3                                 |                 |
|                                   | $H_x$           |
| 5                                 | $H_y$           |
|                                   | H <sub>z</sub>  |

<span id="page-41-2"></span>Table 21: Definition of the <i: field> parameter in the process.dat file.

# 7.2.2 Calculation of frequency spectra: xtransall

If the frequency spectra of points in the TDOM BINARY observer are required, an FFT can be applied to all the time series in impulse.dat and the resulting spectra stored in another binary file called frequency.dat. This is accomplished using the program xtransall, which reads in impulse.dat, the output point information stored in process.dat and the excitation waveform information in excite.dat and then transforms all of the output points:

## \$ xtransall

By default only magnitude data is computed. If phase information is also required then

## \$ xtransall phase

should be used.

The frequency spectra are written to frequency.dat are truncated at a maximum frequency. This maximum frequency is the smallest value of:

- The frequency which corresponds to a mesh size of  $\lambda/5$ .
- Half of the points in the frequency response computed by the FFT.

## 7.2.3 Data extraction at a point or along a line: xtime, xfreq

The programs xtime and xfreq allow the extraction of time series and frequency spectra of a field component, either at a point or along a line, from the impulse.dat and frequency.dat files respectively.

In the process.dat file the parameters n1 and n2 should be set to specify the start and stop time-step or frequency numbers to be extracted. The parameters ilo, ihi, jlo, jhi, klo and khi should be set to the required position(s) and ifield to the required field component as described in Table [21.](#page-41-2) The data is then extracting using

\$ xtime

or

## \$ xfreq phase

if phase information is required.

The output ASCII file written is one of either tdex.dat, tdey.dat or tdez.dat for the time series data or either fdex.dat, fdey.dat or fdez.dat for the frequency spectra. A comment is inserted at the start of the data file showing the indices of the cell(s) of the point(s) chosen. The format of the output file is either time (ns) and field  $(V/m)$  for the time series, or frequency (MHz), logarithmic field (dB V/m or dB  $A/m$ , and field  $(V/m$  or  $A/m$  for the frequency spectra. If phase information is requested another column is written with the phase in degrees. If more than one point is specified, an extra carriage return is inserted after each point, thus allowing a mesh plot of the field along a line against the position along the line and either time or frequency.

## 7.2.4 Data extraction along a line or across plane: xplane, xfplane

The programs xplane and xfplane allow the extraction of the time series and frequency spectra of a field component, either along a line or across a plane, from the impulse.dat and frequency.dat files respectively.

In the process dat file the parameters  $n1$  and  $n2$  should both be set to the time-step or frequency number to be extracted, i.e. both numbers must be the same. The parameters ilo, ihi, jlo, jhi, klo and khi should be set to the required position(s) and ifield to the required field component as described in Table [21.](#page-41-2) The data is then extracting using

\$ xplane

or

```
$ xfplane phase
```
where the phase is needed if phase information is required.

The output ASCII file written is one of either tdex.dat, tdey.dat or tdez.dat for the time series data or either fdex.dat, fdey.dat or fdez.dat for the frequency spectra. A comment is inserted at the start of the data file showing the value of the time (s) or frequency (Hz) chosen. The data is then written in the format of *i*-index, *j*-index, *k*-index and field  $(V/m \text{ or } A/m)$  for both the time series and the frequency spectra. If phase information is requested another column is written with the phase in degrees. If only one coordinate varies, the data will just be along a line.

## 7.2.5 Animated plots: mxplane, mxfplane

The programs mxplane and mxfplane allow the extraction of multiple time series and frequency spectra of a field component, either along a line or across a plane, from the impulse.dat and frequency.dat files respectively.

In the process.dat file the parameters n1 and n2 should be set to specify the start and stop time-step or frequency numbers to be extracted. The parameters ilo, ihi, jlo, jhi, klo and khi should be set to the required position(s) and ifield to the required field component as described in Table [21.](#page-41-2) The data is then extracting using

\$ mxplane

or

## \$ mxfplane phase

if phase information is required.

Each plane of data is written to a separate ASCII file, suitable for plotting as a mesh plot in Gnuplot. The files written are called tds1.dat, tds2.dat, etc, for the time domain data or fds1.dat, fds2.dat, etc, for the frequency domain data. A comment is inserted at the start of each file showing the value of the time (s) or frequency (Hz) chosen. The data is written in the format of the two changing coordinates and the field (V/m or A/m) for both the time series and the frequency spectrum. If only one iteration is specified, that and the next iteration will be output. If only one coordinate varies, the data will just be along a line and will not be suitable for plotting in an animated sequence.

The programs also create a file called plot that contains a Gnuplot script to animate the sequence of data files. This can be loaded into Gnuplot using

## gnuplot> load 'plot'

to view the animation sequence.

## 7.2.6 Interleaving of multiple TDOM BINARY observer data

The tools in the processing toolbox work simultaneously on all the outputs points in the impulse.dat file, regardless of whether they "belong" to different OP directives. Thus, if multiple OP directives of type TDOM BINARY are used in the same simulation and the sets of output cells in the different directives are not disjoint any shared point will be output multiple times. If this is not accounted for then spurious lines can appear in graphs generated from the data. One way to mitigate this problem for lines plots is to sort the data according to position before plotting it.

# 7.3 HDF5 format output

Python and Octave scripts are currently not implemented!

# 8 Tutorial

The mesh files and processing scripts for these examples are included in the Vulture software distribution in the examples directory.

# 8.1 Example 1: Free-space

This first example is worked through in detail.

### 8.1.1 Creating and viewing the mesh

The simplest example of a FDTD simulation is an empty volume of space. Create an ASCII text file called free\_space.mesh with the following contents using a text editor:

```
VM 1.0.0
CE Vulture example: An empty space with absorbing boundaries.
# The mesh extents are 20x40x20 cells.
DM 20 40 20
C<sub>S</sub># Define all external surfaces to be Mur ABCs.
BT XLO MUR
BT XHI MUR
BT YLO MUR
BT YHI MUR
BT ZLO MUR
BT ZHI MUR
# The waveform is a Gaussian pulse with default paramters.
WF wf1 GAUSSIAN_PULSE
# The source is a surface of z-polarised soft electric fields.
EX 0 20 10 10 0 20 source EZ wf1 1.0
# Observe the fields on a plane, every second cell.
OP 0 20 0 40 10 10 plane TDOM_BINARY 2 2 2
# Observe the fields at the centre of the mesh.
OP 10 10 20 20 10 10 centre1 TDOM_ASCII
OP 10 10 20 20 10 10 centre2 FDOM_ASCII
GE
# Run for 400 time-steps.
NT 400
# The mesh size is 1 m.
MS 1.0
EN
```
Now run gvulture

\$ gvulture -p free\_space.mesh

to create Gnuplot format output files and then create an image of the mesh using

```
$ gnuplot mesh.gnp
```
The resulting image in the file mesh.eps is shown in Figure [12.](#page-45-0) The mesh shows the source plane and observer output points. Note the direction arrow on the source plane indicating the direction of polarisation of the electric field.

#### 8.1.2 Running the solver

Now run the solver with the command

\$ vulture -v free\_space.mesh

The progress of the simulation can be checked in the log file vulture.log, though this example should run in only a few seconds.

## 8.1.3 Viewing ASCII data

The time response at the centre of the mesh from the ASCII observer will be stored in the file eh\_centre1\_td.dat. This can be plotted with Gnuplot using

```
set terminal post eps enhanced color 'Helvetica' 18
set output 'free_space_td.eps'
set xlabel 'Time (ns)'
set ylabel 'Electric field, E_z(10,20,10) (V/m)'
plot 'eh_centre1_td.asc' us ($2/1e-9):5 ti '' w l ls 1
```


Vulture example: An empty space with absorbing boundaries.

<span id="page-45-0"></span>Figure 12: Example 1 - Mesh geometry created by gvulture and gnuplot.

creating the image file free space td.eps shown in Figure [13.](#page-46-0) The response exhibits two interesting effects:

- There is a non-physical DC response due to the charging of the grid capacitance associated with the discretised space. FDTD does not include explicit charge dynamics - charges only appear implicitly in the field distributions as determined by Gauss's Law. Hence, since the Gaussian waveform has a DC component it displaces charge across the source plane leaving a net static field on the grid.
- "Spurious" reflections from the edge of the mesh can be seen due to the truncation of the source plane. The source plane only has a finite size and hence the edge launches waves into the mesh.

The frequency spectrum at the centre of the mesh from the ASCII observer will be stored in the file eh centre2 fd.dat. This can be plotted with Gnuplot using

```
set terminal post eps enhanced color 'Helvetica' 18
set output 'free_space_fd.eps'
se xlabel 'Frequency (MHz)'
se ylabel 'Electric field, |E_z|(10,20,10) (dB V/m)'
db20ri( r , i ) = 10.0 * log10 ( r**2 + i**2 )
plot 'eh_centre2_fd.asc' us ($1/1e6):(db20ri($6,$7)) ti '' w l lw 2
```
creating the image file free space fd.eps shown in Figure [14.](#page-46-1) The spurious DC response is now isolated from the physical response. The ripples in the spectrum are again a manifestation of the finite size of the source plane. In order to increase the resolution of the spectrum the simulation needs to be run for more time-steps.

#### 8.1.4 Post-processing binary data in the time-domain

Edit the last line of the process.dat file created by the solver so that the file contains

```
CE Vulture example: An empty space with absorbing boundaries.
1
 0 20 2 0 40 2 10 10 2
1 400
```


<span id="page-46-0"></span>Figure 13: Example 1 - Time response for  $E_z$  at the centre of the mesh.



<span id="page-46-1"></span>Figure 14: Example 1 - Frequency response for  $E_z$  at the centre of the mesh.

0 5.99585e+08 1.49896e+06 1 10 10 20 20 10 10 3

and then extract the time data at the centre of the mesh using

\$ xtime

which should create the output ASCII file tdez.dat. This can be plotted with the previous ASCII data using the Gnuplot script

```
set terminal post eps enhanced color 'Helvetica' 18
set output 'free_space_td2.eps'
set xlabel 'Time (ns)'
set ylabel 'Electric field, E_z(10,20,10) (V/m)'
```
plot 'eh\_centre1\_td.asc' us  $(\$2/1e-9):5$  ti 'ASCII' w l ls 1, \ 'tdez.dat' us 1:2 ti 'Binary' w 1 lw 2

and should give identical results.

#### 8.1.5 Post-processing binary data in the frequency-domain

The frequency response of all the data in the binary impulse.dat file can be calculated using the command

#### \$ xtransall phase

where the phase argument requests that phase information is included - by default xtransall only generates magnitude information. Using the same process.dat as in the previous section the command

#### \$ xfreq phase

will then extract the frequency response at the centre of the mesh into the ASCII file fdez.dat. This can be plotted with

```
se xla 'Frequency (MHz)'
se yla 'Electric field, |E_z|(10,20,10) (dB V/m)'
plot 'fdez.dat' us 1:2 ti '' w l lw 2
se term post eps enhanced color 'Helvetica' 18
se ou 'um_eg1_plotfd.eps'
replot
```
and again should be identical to the output from the ASCII observer.

# 8.2 Example 2: An infinite parallel plate transmission line

Another simple example is a parallel plate transmission line of "infinite" width. The top and bottom plates are represented using PEC boundaries on the external surfaces of the mesh, while the infinite width of the waveguide can by modelled by using PMC boundary conditions on the "sides" of the mesh. Compared to Example 1 we only need to change the boundary conditions on the external surfaces. The complete mesh file is

```
VM 1.0.0
CE Vulture example: An empty parallel-plate waveguide.
# The mesh extents are 20x40x20 cells.
DM 20 40 20
CS# Side walls are PMC.
BT XLO PMC
BT XHI PMC
# End walls are absorbing ABC.
BT YLO PML
BT YHI PML
# Top wals are PEC.
BT ZLO PEC
BT ZHI PEC
# The waveform is a Gaussian pulse with default paramters.
WF wf1 GAUSSIAN_PULSE
# The source is a surface of z-polarised soft electric fields.
EX 0 20 10 10 0 20 source EZ wf1 1.0
# Observe the fields on a plane, every second cell.
OP 0 20 0 40 10 10 plane TDOM_BINARY 2 2 2
# Observe the fields at the centre of the mesh.
OP 10 10 20 20 10 10 centre1 TDOM_ASCII
OP 10 10 20 20 10 10 centre2 FDOM_ASCII
GE
# Run for 400 time-steps.
NT 400
```
### # The mesh size is 1 m. MS 1.0 EN

and the output from gvulture looks exactly like the free-space example is Figure [12.](#page-45-0) We have also chosen to use PML absorbing boundaries on the open ends of the waveguide rather than the Mur ABCs used in Example 1.





<span id="page-48-0"></span>Figure 15: Example 2 - Time response for  $E_z$  at the centre of the mesh.

Figure 16: Example 2 - Frequency response for  $E_z$  at the centre of the mesh.

The processing steps are exactly the same as for Example 1. The time response at the centre of the mesh is shown in Figure [15](#page-48-0) and the frequency spectrum at the smae point in Figure [15.](#page-48-0) The "spurious" reflections from the edge of the mesh in the previous example have now disappeared since the source plane is now consistent with the boundary conditions on the external surfaces of the mesh. The DC response has also disappeared since the charge on the source plane edges can discharge into the PEC plates. A soft electric field source can therefore excite an accutate TEM wave in a parallel-plate waveguide structure,

however, it launches a wave in both directions!



<span id="page-49-0"></span>Figure 17: Example 2 - Time response of  $E_z$  across the  $k = 10$  plane at time-step 60.

The time response of the electric field can be extract from the binary observer over a whole plane. To accomplish this for time-step number 60 modify the process.dat file to contain

```
CE Vulture example: An empty parallel-plate waveguide.
1
0 20 2 0 40 2 10 10 2
60 60
0 5.99585e+08 1.49896e+06
1
0 20 0 40 10 10 3
```
Here the parameters n1 and n2 are used to set the required output time-step and the bounding box in the last line ranges over the whole observer plane. Now run the xplane program

### \$ xplane

This should generate an output file called tdez.dat containing mesh plot data for the field that can be plotted using the Gnuplot script

```
set terminal post eps enhanced color 'Helvetica' 18
set output 'parallel_plate_tds60.eps'
set xlabel 'i (cells)'
set ylabel 'j (cells)'
set zlabel 'Electric field, E_z(i,j,10) (V/m)' rotate by 90 offset -1.5,1.5
set ticslevel 0
set title "Time-step: 60"
set view 120,60
splot 'tdez.dat' us 1:2:4 ti '' w l ls 1
```
The result is shown in Figure [17.](#page-49-0) Note how the soft field source excites a wave of amplitude one in both directions normal to the surface over which it is imposed. The amplitude depends on the CFLN as can be investigated by adding a CN directive, such as

#### CN 0.99

to section 3 of the mesh file. As the CFLN is increased the amplitude of the TEM wave launched is reduce and visa-vera.

The presence of both forward and backwards waves may or may not be a problem depending on the application. The use of a PML in the open ends of the waveguide is advisable to provide low reflection of the backward travelling wave into the mesh.

## 8.3 Example 3: An accurate plane-wave in free-space

To launch an accurate plane-wave in free-space a total-field scattered-field plane-wave source can be used. Starting from the mesh of Example 1 again we replace the distributed electric field source with a plane-wave source and also change the external surfaces to be PMLs rather than Mur ABCs. The complete mesh is now

```
VM 1.0.0
CE Vulture example: Plane wave in free-space.
# The mesh extents are 20x40x20 cells.
DM 20 40 20
C<sub>S</sub># Define all external surfaces to be PML.
BT XLO PML
BT XHI PML
BT YLO PML
BT YHI PML
BT ZLO PML
BT ZHI PML
# The waveform is a Gaussian pulse with default paramters.
WF wf1 GAUSSIAN_PULSE
# Plane-wave in +ve y direction, Ez polarisation.
PW 3 17 3 37 3 17 pw1 wf1 90 90 90 111111 1.0 0.0
# Observe the fields on a plane, every cell.
OP 0 20 0 40 10 10 plane TDOM_BINARY 1 1 1
# Observe the fields at the centre of the mesh.
OP 10 10 20 20 10 10 centre1 TDOM_ASCII
OP 10 10 20 20 10 10 centre2 FDOM_ASCII
GE
# Run for 400 time-steps.
NT 400
# The mesh size is 1 m.
MS 1.0
EN
```
The mesh now looks like Figure [18.](#page-51-0) The TFSF surface is shown as a box on the mesh with arrows at the corner where the plane-wave first enters the total-field zone depicting the direction of propagation (open arrow), polarisation direction the electric field (closed arrow) and polarisation direction of the magnetic field (double open arrow). A higher density of observation points have been requested by changing the stride lengths in the binary observer to one cell.

Proceeding with the same proccessing steps as before we obtain the time response and frequency spectrum at the centre of the mesh shown in Figures [19](#page-52-0) and [20](#page-52-1) respectively. Compared to Example 1 we observe that a substantially improved plane-wave is excited within the total-field zone in the mesh. Up to a frequency corresponding to a mesh-size of one-tenth of a wavelength the amplitude of the wave remains within 0.1 dB of the ideal value.

Surface plots of the electric field over the plane of the binary observer can be obtained as before. For example to get the response at time-step 60 modify the process.dat file to be

```
CE Vulture example: Plane wave in free-space.
1
0 20 1 0 40 1 10 10 1
60 60
0 5.99585e+08 1.49896e+06
1
 0 20 0 40 10 10 3
```


Vulture example: Plane wave in free-space.

<span id="page-51-0"></span>Figure 18: Example 3 - Mesh geometry created by gvulture and gnuplot.

and run

#### \$ xplane

Figure [21](#page-53-0) shows surface plots for time-steps 60 and 100. Notice how there is no field in the scattered field zone since there are no scattering objects inside the TFSF surface. The plane-wave is completely "reabsorbed" by the TFSF surface at  $j = 37$ .

An animation of the whole time response in the plane can be created by changing the time-step intervals in the process.dat file to cover the whole time range:

```
CE Vulture example: Plane wave in free-space.
1
0 20 1 0 40 1 10 10 1
1 400
0 5.99585e+08 1.49896e+06
1
0 20 0 40 10 10 3
```
Using the command

#### \$ mxplane

a series of 400 data files with names tds<nnn>.dat are created that can be viewed in sequence to create an animation.

\$ gnuplot gnuplot> load 'plot'

## 8.4 Example 4: Transmission and reflection from a dielectric slab

This example calculates the reflection and transmission coefficient of an infinite dielectric slab to a normally incident TEM wave. To simulate the infinite extent of the slab it is placed in the cross-section of a parallelplate waveguide as was used in Example 2. Since the problem is invariant along the directions parallel to the plane of the slab only one cell needs to be used in those directions. The effectively one-dimensional mesh is given by

VM 1.0.0 CE Vulture example: Parallel plate waveguide with dielectric slab



<span id="page-52-0"></span>Figure 19: Example 3 - Time response for  $E_z$  at the centre of the mesh.



<span id="page-52-1"></span>Figure 20: Example 3 - Frequency response for  $E_z$  at the centre of the mesh.

DM 200 1 1 GS # Parallel-plate waveguide boundary conditions. BT XLO PML BT XHI PML BT YLO PEC BT YHI PEC BT ZLO PMC BT ZHI PMC # Simple dielectric medium. MT dielectric SIMPLE 2.0 0.01 1.0 MB 50 150 0 1 0 1 dielectric # Gaussian source.



<span id="page-53-0"></span>Figure 21: Example 3 - Time responses of  $E_z$  across the  $k = 10$  plane at time-steps 60 and 100.



Vulture example: Parallel plate waveguide with dielectric slab

<span id="page-53-1"></span>Figure 22: Example 4 - Mesh geometry created by gvulture and gnuplot.

```
WF wf1 GAUSSIAN_PULSE
# Partial TF/SF.
PW 30 170 0 1 0 1 pw1 wf1 90 0 180 110000 1.0 0.0
# Observe field reflected field if SF zone.
OP 20 20 0 0 0 0 ref1 TDOM_ASCII
OP 20 20 0 0 0 0 ref2 FDOM_ASCII
# Observe transmitted field in TF zone.
OP 160 160 0 0 0 0 trans1 TDOM_ASCII
OP 160 160 0 0 0 0 trans2 FDOM_ASCII
# Observe field along waveguide.
OP 0 200 0 0 0 0 line TDOM_BINARY 1 1 1
GE
NT 5000
MS 0.01
EN
```
The lossy dielectric has a relative permittivity of 2 and a conductivity of 10 mS/m. These parameters are defined by the MT directive. The material parameters are applied to the mesh between  $i = 50$  and  $i = 150$  by the MB directive. A TEM wave could be excited using a plane of soft electric field sources as was done in Example 2. However, in this case we use a partial TFSF plane-wave source, which has the advantage of allowing the reflected wave to be observed directly in the scattered field zone.

The mesh geometry is shown in Figure [22.](#page-53-1) Only the XLO and XHI surfaces of the TFSF are active, as set by the mask 110000 in the PW directive. The wave is excited at  $i = 30$  and the upper surface of the TF-SF boundary is at  $i = 170$ . Therefore the transmitted wave can be observed at the point  $i = 160$  in the total field zone and the reflected wave at  $i = 20$  in the scattered field zone.



<span id="page-54-0"></span>Figure 23: Example 4 - Time response at time-steps 50, 100, 150 and 350.

The solver is run in the usual way and the response along the length of the waveguide from the binary observer is extracted at all time steps by editing the process.dat to read

```
CE Vulture example: Parallel plate waveguide with dielectric slab
1
 0 200 1 0 0 1 0 0 1
1 1000
```


<span id="page-55-0"></span>Figure 24: Example 4 - Scattering parameters of the dielctric slab: Magntitude (top) and phase (bottom). The phase reponse is only shown to 2 GHz for clarity.

0 5.99585e+10 1.19917e+07 0.01 0 200 0 0 0 0 2

and then running

#### \$ mxplane

to obtain data files tds<nnn>.dat for the response along a line at each time-step. These can be plotted using, for example, the Gnuplot script

```
set terminal post eps enhanced color 'Helvetica' 16
set output 'frame100.eps'
set style data lines
set xrange [0:200]
set yrange [-1:1]
```

```
set xlabel 'x (cells)'
set ylabel 'E_y (V/m)'
set title 'time-step: 100'
set arrow 1 from 30,-1 to 30,1 nohead lt 0
set arrow 2 from 170,-1 to 170,1 nohead lt 0
set label 1 'SF' at 26,-0.9 left rotate by 90
set label 2 'TF' at 34,-0.9 left rotate by 90
set label 3 'TF' at 166,-0.9 left rotate by 90
set label 4 'SF' at 174,-0.9 left rotate by 90
set arrow 3 from 50,-1 to 50,1 nohead lt 0
set arrow 4 from 150,-1 to 150,1 nohead lt 0
set label 5 'Free-space' at 46,-0.9 left rotate by 90
set label 6 'Dielectric' at 54,-0.9 left rotate by 90
set label 7 'Dielectric' at 146,-0.9 left rotate by 90
set label 8 'Free-space' at 154,-0.9 left rotate by 90
plot 'tds100.dat' us 1:2 ti ''
```
Figure [23](#page-54-0) shows the spatial variation of the field at a series of time-steps. At time-step 50 the pulse is emerging from the lower TFSF boundary and is approaching the dielectric's surface. By time-step 100 the pulse is approximately centred on the lower dielectric interface and the reflected pulse is just becoming visible. At time-step 150 the reflected pulse has propagated away into the lower scattered-field zone and the transmitted pulse has cleared the dielectric interface and propagates into the medium. The attenuation of the pulse due to the dielectric loss can be observed at time-steps 200 and 250. At time-step 350 the incident field has entered the upper scattered-field zone. Note, if there were no dielectric present this field would be cancelled by the field which propagates through the total-field zone, giving zero scattered field. However, the dielectric slab attenuates and delays the wave in the total-field zone and so a non-zero scattered field is manifested in the upper scattered-field zone.

The scattering parameters of the slab with reference to its surfaces are:

$$
S_{11} = \frac{E_z^-(50)}{E_z^+(50)} \tag{8.1}
$$

$$
S_{21} = \frac{E_z^+(150)}{E_z^+(50)}.\tag{8.2}
$$

These can be calculated analytically [\[11\]](#page-66-5). The fields at the slab surfaces are related to those at the observation and excitation points by

$$
\frac{E_z^+(160)}{E_z^+(150)} = e^{-j\omega \cdot 10 \Delta t/c_0}
$$
\n(8.3)

$$
\frac{E_z^+(50)}{E_z^+(29)} = e^{-j\omega \cdot 21\Delta t/c_0}
$$
\n(8.4)

$$
\frac{E_z^-(20)}{E_z^-(50)} = e^{-j\omega \cdot 30 \Delta t/c_0}
$$
\n(8.5)

Note that in terms of the phase reference the plane-wave is effectively launched from one cell inside the scattered field zone, hence the field  $E_z^+(29)$  is used as the reference for the FDTD spectra. It is usually more reliable to calibrate the incident field at the location of the front face of the dielectric by running a reference simulation with the dielectric slab removed and an observation point at  $i = 50$  rather than relying on knowledge on the internal implementation of the code. Using the above we find that

$$
S_{11} = e^{j\omega \cdot 51 \Delta t/c_0} \frac{E_r^-(20)}{E_z^+(29)}
$$
(8.6)

$$
S_{21} = e^{\jmath \omega \cdot 31 \Delta t/c_0} \frac{E_z^+(160)}{E_z^+(29)}.
$$
\n(8.7)

Figure [24](#page-55-0) shows the scattering parmeters of the dielectric slab compared with the analytical solution. It can be seen that the lower frequencies give accurate results whereas the higher frequency results are not so reliable. For high accuracy simulations the standard  $\lambda/10$  rule of thumb is not sufficient and a maximum mesh edge length of  $\lambda/20$  should not be exceeded at the highest frequency of interest. In the analytic result

a slab thickness of 101 cells was used - one more than the thickness of the FDTD bounding box. This is because the simulation was run using an executable without the averaged media support enabled and hence the dielectric material parameters are applied "in-full" on both the front and back (XLO and XHI) surfaces of the slab. Applied in this way the material parameters are effective to about half a cell from these surfaces leading to an overall effective thickness one cell greater than the bounding box thickness.

# 8.5 Example 5: Transmission and reflection from a thin boundary

## 8.6 Example 6: Radiation from an infinitesimal Hertzian dipole

8.7 Example 7: Radiation from a half-wave dipole

## 8.8 Example 8: Plane-wave penetration through an aperture

This example looks at the transmission through a small square aperture in an infinite PEC plate using a truncated plane-wave excitation. The mesh for the problem is:

```
VM 1.0.0
CE Vulture example: Aperture in infinite plane
DM 60 40 40
GS
# Free-space boundaries.
BT XLO PML
BT XHI PML
BT YLO PML
BT YHI PML
BT ZLO PML
BT ZHI PML
# Solid PEC sheet
#TB 30 30 0 40 0 40 PEC
# PEC sheet with 2x2 cell square aperure.
# TB 30 30 0 40 0 19 PEC
# TB 30 30 0 40 21 40 PEC
# TB 30 30 0 19 19 21 PEC
# TB 30 30 21 40 19 21 PEC
# Differentiated Gaussian pulse waveform
WF wf1 DIFF_GAUSSIAN_PULSE 1.0
# Full plane-wave for empty mesh case.
PW 5 55 5 35 5 35 pw1 wf1 90.0 0.0 90.0 111111 1.0 0.0
# Truncated incident wave.
PW 5 30 5 35 5 35 pw1 wf1 90.0 0.0 90.0 101111 1.0 0.0
# Add reflected wave using second plane-wave, delay by flight time along bbox.
# PW 5 30 5 35 5 35 pw2 wf1 90.0 180.0 270.0 101111 1.0 8.339e-11
# Magnetic moment source using differential waveform
# dt = 1.667818e-12
# dl = 1e-3
# a = 2 * d1# alpha = 4 * a^3 / 3 / pi^1.5
# Hsc = 2 / eta0# m = -alpha * Hsc# sigma = 5 * sqrt(2) * dt# Imdl = mu0 * m * 1 / sigma * d/dt WF
# delay = (25 + 0.5) * d1 / c0# WF wf2 RICKER_WAVELET 1.0
# EX 30 31 19 19 19 21 dipmom IMDY wf2 1.0836e-06 8.5059e-11
# Observe incident field point.
OP 29 29 20 20 20 20 inc1 TDOM_ASCII
OP 29 29 20 20 20 20 inc2 FDOM_ASCII
# Observe transmitted field on axis of aperture.
OP 50 50 20 20 20 20 trans1 TDOM_ASCII
```
OP 50 50 20 20 20 20 trans2 FDOM\_ASCII # Observe fields in plane through aperture. OP 0 60 0 40 20 20 plane TDOM\_BINARY 1 1 1 GE NT 200 MS 0.001 EN

Some parts of the mesh are commented out and will be used later. The active parts currently simulate a plane-wave in free space to use as a reference case. The simulation is run and the data post-processed using the same steps as in the previous examples. The tangential magnetic field spectrum at mesh point  $(29, 20, 20)$  is shown in Figure [26](#page-59-0) and as expected it has a uniform magnitude of  $1/\eta_0$  A/m.



#### Vulture example: Aperture in infinite plane

<span id="page-58-0"></span>Figure 25: Example 8 - Geometry of the aperture transmission problem, showing the TFSF bounding box, aperture and observation points either side of the aperture plane. The plane observer points are not shown for clarity.

Now the problem space is cut by introducing a solid PEC surface in the  $i = 30$  plane

```
# Solid PEC sheet
TB 30 30 0 40 0 40 PEC
...
```
and replacing the full TFSF surface by a partial surface

```
...
# Truncated incident wave.
PW 5 30 5 35 5 35 pw1 wf1 90.0 0.0 90.0 101111 1.0 0.0
...
```
which only extends as far as the PEC plate and where the XHI surface has been deactivated using the mask 101111. The wave at time-step 120, after the wave has reflected from the PEC surface, is shown in the top-left of Figure [26.](#page-59-0) It can be seen that it is quite distorted compared to the expected superposition of incident and reflected plane-waves. This is because the TFSF surface only injects the incident field into the total field region and the reflected field "spills out" into the scattered field region. The effect can also be seen in the amplitude spectrum of the tangential magnetic field at the PEC surface, which exhibits some ripples compared to the expected flat value of  $2/\eta_0$ . This can be thought of as arising from secondary Huygen's sources located along the edges where the TFSF and PEC plate intersect; on these edge the field injected



<span id="page-59-0"></span>Figure 26: Example 8 - Spatial response of electric across a plane for a plane-wave incident on a solid PEC plate at time step 120 without (top left) and with (top right) the reflected wave included in the TF-SF exciation. The frequency reponse of the tangential magnetic field directly next to the PEC plane is shown below.

by the TFSF surface does not satisfy the PEC boundary conditions. To remedy this the reflected wave can be included in the field injected by the TFSF surface. In Vulture this can be accomplished by adding a second PW directive with a wave travelling in the opposite direction, with the opposite polarisation:

... # Add reflected wave using second plane-wave, delay by flight time along bbox. PW 5 30 5 35 5 35 pw2 wf1 90.0 180.0 270.0 101111 1.0 8.339e-11 ...

The wave is also delayed by a time of  $(30-5)\Delta/c_0 = 8.339$  ps, corresponding to the propagation time along the partial Huygen's surface, so the reflected wave begins as the incident wave strikes the PEC plate. The wave at time-step 120 with this reflected wave included is shown in the top-right of Figure [26.](#page-59-0) This shows a substantial improvement in the wave shape. There is still some distortion due to the fact that the incident wave undergoes dispersion as it propagates towards the PEC plate that is not accounted for by the injected reflected wave. In a "real" simulation the length of the TFSF box would be reduced to one or two cells as in

...

# Add reflected wave using second plane-wave, delay by flight time along bbox. PW 28 30 5 35 5 35 pw2 wf1 90.0 180.0 270.0 101111 1.0 8.339e-11 # Truncated incident wave. PW 28 30 5 35 5 35 pw1 wf1 90.0 0.0 90.0 101111 1.0 0.0 ...

in order to mitigate these dispersion effect. Here, for the sake of presentation, the TFSF has been made sufficiently long to allow the incident and reflected waves to be seen clearly.

Now that we have an accurate incident field, a small square aperture of side length  $a = 2\Delta l$  is cut in the PEC plate by replacing the solid plate TB directive with

... # PEC sheet with 2x2 cell square aperure. TB 30 30 0 40 0 19 PEC TB 30 30 0 40 21 40 PEC TB 30 30 0 19 19 21 PEC TB 30 30 21 40 19 21 PEC ...

The mesh for this case is shown in Figure [25.](#page-58-0) The transmitted wave is shown in the top-left of Figure [27](#page-61-0) at time-step 130. A "spherical" wave can be seen radiating from the aperture. The transmitted electric field is observed at mesh cell (50, 20, 20) and its time series is shown in the graph in the lower part of Figure [27.](#page-61-0)

The field radiated by a small aperture can be approximated by the field from a set of equivalent dipole moments [\[12\]](#page-66-6). For a normally incident plane-wave only the magnetic dipole moment polarised in the direction of the magnetic field is excited and this is given by

$$
m_y = \alpha_{\text{M};yy} H_y^{\text{SC}} = 2\alpha_{\text{M};yy} H_y^{\text{inc}},\tag{8.8}
$$

where  $\alpha_{M;yy}$  is the magnetic polarisability of the aperture and  $H_y^{\text{SC}}$  is the short-circuited magnetic field on the incident field side, equal to twice the incident magnetic field  $H_y^{\text{inc}}$  [\[13\]](#page-66-7). The polarisability of the aperture is approximately

$$
\alpha_{\mathrm{M};yy} = \frac{4a^3}{3\sqrt{\pi}^3},\tag{8.9}
$$

where  $\alpha$  is the side length. The current moment of the equivalent infinitesimal magnetic dipole is given by

$$
I_{\mathcal{M};y}(t) dy = \mu_0 \frac{dm_y}{dt}
$$
\n(8.10)

and from the above we can write this

$$
I_{\mathcal{M};y}(t) \, \mathrm{d}y = \mu_0 \frac{8a^3}{3\sqrt{\pi}^3} \frac{\mathrm{d}H_y^{\text{inc}}}{\mathrm{d}t}.\tag{8.11}
$$

In the "real" aperture similation the incident waveform was a differentiated Gaussian pulse,  $\psi_{\text{DGP}}(t)$ , so accounting for the delay between the waveform and the location of the equivalent dipole moment of the aperture is

$$
I_{\mathcal{M};y}(t) \, \mathrm{d}y = \left. \frac{\mu_0}{\eta_0} \frac{8a^3}{3\sqrt{\pi}^3} \, \frac{\mathrm{d}\psi_{\text{DGP}}}{\mathrm{d}t} \right|_{t=25.5\Delta t} . \tag{8.12}
$$

The derivative of the differentiated Gaussian pulse waveform is just a scaled version of the Ricker Wavelet waveform so finally

$$
I_{\mathrm{M};y}(t) \, \mathrm{d}y = \frac{\mu_0}{\eta_0} \frac{8a^3}{3\sqrt{\pi}^3} \frac{1}{\sigma_{\mathrm{GP}}} \psi_{\mathrm{RW}}(t - 25.5\Delta t). \tag{8.13}
$$

This can be implemented in the simulation by removing the plane-wave source, closing the aperture in the PEC plate and then introducing the magnetic moment source

... # WF wf2 RICKER\_WAVELET 1.0 EX 30 31 19 19 19 21 dipmom IMDY wf2 1.0836e-06 8.5059e-11 ...



<span id="page-61-0"></span>Figure 27: Example 8 - Spatial response of the electric field transmitted by the aperture across a plane at time-step 130 for the physical aperture (top left) and equivlanet dipole (top right). The time reponse of the electric field at a point on the axis of the aperture is shown below.

The top-right part of Figure [27](#page-61-0) shows the wave radiated by this equivalent dipole moment source. While it is similar to that of the aperture it is not the same. Looking at the wave at (50, 20, 20) is can be seen that the wave from the equivalent dipole moment is larger and slightly ahead of that of the physical aperture simulation. The reasons for this are:

- The effective size of the aperture on the discretised FDTD grid is smaller than  $a = 2\Delta l$ . The PEC conditions are enforced on the aperture edges, but have an effective "range of influnence" of the order of  $\Delta l/2$ . The effective size of the aperture is there somewhat smaller than  $2\Delta l$ .
- The grid dispersion of the incident wave in the physical aperture simulation is not accounted for in the equivalent dipole simulation.
- The physical aperture may have some dispersion (delay) associated with its finite size.

Good agreement between the physical aperture simulation and equivalent dipole simulation can be obtained using an effective aperture size of  $a = 1.49\Delta l$  and adding an excess delay of 7 ps, as shown in Figure [27.](#page-61-0) By varying the number of cells in the aperture while keeping the physical aperture size fixed it can be shown that at least 8 cells per side are required to obtain an accuracy of 12 % or better.

# $\bigcup^Z$  x

## 8.9 Example 9: Shielding effectiveness of an enclosure with an aperture

<span id="page-62-0"></span>Figure 28: Example 9 - Mesh geometry created by gvulture and gmsh.

This example consists of a metal sided enclosure of dimensions  $300 \times 120 \times 300$  mm with a  $200 \times 30$  mm slot in the centre of one side. The enclosure is illuminated by a plane-wave which is normally incident on the slot. The input mesh file is:

```
VM 1.0.0
CE Vulture example: Enclosure with slot.
DM 80 100 44
GS
# Front face of enlcosure with slot..
TB 10 70 30 30 10 19 PEC
TB 10 70 30 30 25 34 PEC
TB 10 20 30 30 19 25 PEC
TB 60 70 30 30 19 25 PEC
# Back face.
TB 10 70 90 90 10 34 PEC
# Top face.
TB 10 70 30 90 34 34 PEC
# Bottom face.
TB 10 70 30 90 10 10 PEC
# Left face.
TB 10 10 30 90 10 34 PEC
# Right face.
TB 70 70 30 90 10 34 PEC
# Gaussian waveform.
WF wf1 GAUSSIAN_PULSE
# Incident plane-wave.
PW 15 65 10 30 14 30 pwinc wf1 90.0 90.0 90.0 111011 1.0 0.0
# Reflected plane-wave.
PW 15 65 10 30 14 30 pwref wf1 90.0 270.0 270.0 111011 1.0 3.3356e-10
# Observers at centre of top face.
OP 40 40 60 60 22 22 centre1 TDOM_ASCII
OP 40 40 60 60 22 22 centre2 FDOM_ASCII
# Observer for field in the slot.
OP 40 40 29 29 22 22 slot1 TDOM_ASCII
OP 40 40 29 29 22 22 slot2 FDOM_ASCII
GE
NT 80000
```
#### MS 5e-3 EN

The enclosure is constructed from PEC surface materials. A partial TFSF is again used to introduce the plane-wave excitation of the slot and the reflected wave and we observe the fields in the slot and at the centre of the enclosure. The geometry as viewed in Gmsh is shown in Figure [28.](#page-62-0) The Gmsh input file can be created using

\$ gvulture -e -p -m enclsoure\_se.mesh

and then viewed in Gsmh with

#### \$ gmsh mesh.msh

The time response of the observation point at the centre of the enclosure is shown in Figure [29.](#page-64-0) The attenuation of the signal is entirely due to re-radiation of energy back through the slot since there are no other loss mechanisms inside the enclosure. The shielding effectiveness of the enclosure can be defined as the amount by which the incident field is attenuated by the enclosure's presence. If the incident plane-wave field amplitude is  $|E^{inc}|$  and the internal field amplitude is  $|E^{int}|$  then we can define

$$
S_{\rm E} = \frac{|E^{\rm inc}|}{|E^{\rm int}|}. \tag{8.14}
$$

In order to obtain an accurate frequency spectrum from a time response a rule-of-thumb is that the time response should have decayed by 60 dB from its maximum value. The response in Figure [29](#page-64-0) has only decayed by about 35 dB so a spectrum obtained from it should be treated with caution.

Clearly the field inside the enclosure is highly non-uniform so the shielding effectiveness depends on the position within the enclosure. Figure [30](#page-64-1) shows the shielding effectiveness at the centre of the cavity, which since the incident field amplitude is unity is just the reciprocal of  $|E_z(40, 60, 22)|$ , compared to the prediction of a simple analytical model [\[14\]](#page-66-8). The two results agree well up to just over 1 GHz. Beyond this frequency the analytic model becomes invalid due to the excitation of higher order modes.

Once a cavity is loaded by contents the response will change. To respresent this effect we introduce a 100 mm sided cube of radio absorbing material into the enclsoure. The material is a commercially available absorber, Eccosorb LS22 [\[15\]](#page-66-9). The material's complex permittivity, obtained from the manufacturer's datasheet, is shown in figure [31](#page-65-0) together the result of applying a genetic algorithm to fit the complex permittivity to a third order Debye dispersion relationship. The Debye parameters obtained are given in Table [22.](#page-63-0)

| $\kappa$ | $\Delta \epsilon_k$ | $\tau_k$               | $r_k$           | $p_k$            |
|----------|---------------------|------------------------|-----------------|------------------|
| Ξ.       | $-1$                |                        | $(rads^{-1})$   | $(rads^{-1})$    |
|          | 0.00109             | $0.1$ s                | 5.45118e-03     | $-1.00000e + 01$ |
| 2        | 6.632               | $0.1062$ ns            | $3.12249e+10$   | $-9.41658e+09$   |
| 3        | 3.27                | $15.781 \,\mathrm{ps}$ | $1.03602e + 11$ | $-6.33664e+10$   |

<span id="page-63-0"></span>Table 22: Example 9 - Debye model for LS22 RAM:  $\epsilon_{\infty} = 1, \sigma = 0.344191$ .

The absorber is introduced to the mesh by adding the following two directives

```
MT ls22 DEBYE "LS22-3pole.prm"
MB 30 50 60 80 12 32 ls22
```
The Debye parameters are then entered into an ASCII file called LS22-3pole.prm which should contain:

```
3 1.00000e+00 3.44191e-01 1.00000e+00
1.03602e+11 0.00000e+00 -6.33664e+10 0.00000e+00
5.45118e-03 0.00000e+00 -1.00000e+01 0.00000e+00
3.12249e+10 0.00000e+00 -9.41658e+09 0.00000e+00
```
Running the simulation again we obtain the results denoted by "Loaded" in Figures [29](#page-64-0) and [30.](#page-64-1) The time response now decays much more quickly than that of the empty enclosure and the shielding effectiveness at the centre of the enclosure is generally increased by 20 dB. The resonant features are smeared out and damped by the absorber, though some sharp features remain and there are frequencies at which the shielding has decreased.

TBC:



Figure 29: Example 9 - Time response at centre of enclosure. The amplitude of the electric field is plotted in decibels, showing the long exponential decay of the field inside the unloaded enclosure compared to the rapid decay when the enclosure is loaded with an LS22 block.

<span id="page-64-0"></span>

<span id="page-64-1"></span>Figure 30: Example 9 - Shielding effectiveness at centre of enclosure with and without loading. The analytic model is only valid up to the first resonance of the enclosure.

- Decay time calculation and graph annotation.
- Required energy decay for reliable DFT.
- Partial Q-factor and decay rate of aperture.
- Partial Q-factor and decay rate of LS22.



<span id="page-65-0"></span>Figure 31: Example 9 - Complex permittivity of LS22 RAM comparing manufacturer's data to a third order Debye model obtained using a genetic algorithm.



<span id="page-65-1"></span>Figure 32: Cross-section of the shielded microstrip line.

# 8.10 Example 10: Stripline and microstrip line parameters

This example considers a shielded microstrip structure, as used by Gedney [\[16\]](#page-66-10), and shown in Figure [32.](#page-65-1) The dimension of the structure are  $w = 0.254$  mm,  $h = 0.254$  mm,  $h_a = 0.762$  mm,  $s = 2.11$  mm and the permittivity is  $\epsilon_r = 2.2$ . The structure is modelled on a grid with  $\Delta l = 0.042333$  mm giving an overall grid of dimensions  $50 \times 24 \times 400$  and time step of  $\Delta t = 70.5555$  fs. The line was excited using a plane of y-polarised electric field points underneath the track at the centre of the mesh in the  $z$  direction. A broad Gaussian pulse was used, with a bandwidth of 450 GHz. PML boundaries were used to terminate the problem space at the lower and upper z boundaries. This structure therefore allows the near normal incidence reflection the inhomogeneous PML to be studied.

Replacing the PECs on the sides of the mesh with PML we obtain the open microstrip structure shown in Figure [33.](#page-66-11)

[TBD: Determination of voltage, current, characterstic impedance, propagation constant and effective permittivity.]

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<span id="page-66-11"></span>Figure 33: Cross-section of the microstrip line.

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