

HVDC

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Chapter 1

File Index

1.1 File List

Here is a list of all files with brief descriptions:

test_3.h	Test case with oil and paper layers and oil filled cubic butt gaps	3
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Chapter 2

File Documentation

2.1 test_3.h File Reference

Test case with oil and paper layers and oil filled cubic butt gaps.

```
#include <tmesh_3d.h>
#include <simple_connectivity_3d_thin.h>
```

Functions

- static int [uniform_refinement](#) (tmesh_3d::quadrant_iterator q)
Uniform refinement function.
- static int [refinement](#) (tmesh_3d::quadrant_iterator quadrant)
Performs local refinement at the interfaces.
- static int [coarsening](#) (tmesh_3d::quadrant_iterator quadrant)
Performs local coarsening, leaves interfaces refined.
- double [epsilon_fun](#) (const double &x, const double &y, const double &z)
Return the value of epsilon at the point (x,y,z)
- double [sigma_fun](#) (const double &x, const double &y, const double &z)
Return the value of sigma at the point (x,y,z)
- std::vector< size_t > [find_idx](#) (tmesh_3d &tmsh, std::vector< std::vector< double > > &points, std::vector< std::vector< double > > &tols, const size_t &N_rhos)
Find the global index of the points given by the vector 'points'.

Variables

- constexpr int [NUM_REFINEMENTS](#) = 5
Level for global refinement.
- constexpr int [maxlevel](#) = 6
Level for local refinement.
- constexpr int [minlevel](#) = 4
Level for global coarsening.
- constexpr double [DELTAT](#) = 50.0
Temporal time step.

- constexpr double `T` = 5000
Final time of simulation.
- constexpr double `tau` = 50.0
Time constant for boundary conditions.
- constexpr bool `save_sol` = true
If set to 'true' saves data for Paraview visualization.
- constexpr double `epsilon_0` = 8.8542e-12
Permittivity vacuum.
- constexpr double `epsilon_r_1` = 2.0
Permittivity oil.
- constexpr double `epsilon_r_2` = 4.0
Permittivity paper.
- constexpr double `sigma` = 3.21e-14
Conductivity.
- constexpr double `z_oil` = 5e-5
Thickness oil layer.
- constexpr double `z_paper` = 3e-4
Thickness paper layer and butt gaps side length.
- constexpr double `tol` = 1e-5
Tolerance value for refinement and point selection.
- constexpr size_t `N_rhos` = 6
Number of points to select for output.
- std::vector< size_t > `rho_idx`
- std::vector< std::vector< double > > `points` {{5e-4,5e-4,0.0},{5e-4,5e-4,`z_paper`},{5e-4,5e-4,`z_oil`+`z_paper`},{5e-4,5e-4,`z_oil`+2*`z_paper`},{5e-4,5e-4,2*`z_oil`+2*`z_paper`},{5e-4,5e-4,1e-3}}
- std::vector< std::vector< double > > `tols` {{1e-4,1e-4,`tol`},{1e-4,1e-4,`tol`},{1e-4,1e-4,`tol`},{1e-4,1e-4,`tol`},{1e-4,1e-4,`tol`},{1e-4,1e-4,`tol`}}
- bool `extra_refinement` = true
true for all cases except Test 1

2.1.1 Detailed Description

Test case with oil and paper layers and oil filled cubic butt gaps.

2.1.2 Function Documentation

2.1.2.1 coarsening()

```
static int coarsening (
    tmesh_3d::quadrant_iterator quadrant ) [static]
```

Performs local coarsening, leaves interfaces refined.

2.1.2.2 epsilon_fun()

```
double epsilon_fun (
    const double & x,
    const double & y,
    const double & z )
```

Return the value of epsilon at the point (x,y,z)

2.1.2.3 find_idx()

```
std::vector< size_t > find_idx (
    tmesh_3d & tmsh,
    std::vector< std::vector< double > > & points,
    std::vector< std::vector< double > > & tols,
    const size_t & N_rhos )
```

Find the global index of the points given by the vector 'points'.

2.1.2.4 refinement()

```
static int refinement (
    tmesh_3d::quadrant_iterator quadrant ) [static]
```

Performs local refinement at the interfaces.

2.1.2.5 sigma_fun()

```
double sigma_fun (
    const double & x,
    const double & y,
    const double & z )
```

Return the value of sigma at the point (x,y,z)

2.1.2.6 uniform_refinement()

```
static int uniform_refinement (
    tmesh_3d::quadrant_iterator q ) [static]
```

Uniform refinement function.

2.1.3 Variable Documentation

2.1.3.1 DELTAT

```
constexpr double DELTAT = 50.0 [constexpr]
```

Temporal time step.

2.1.3.2 epsilon_0

```
constexpr double epsilon_0 = 8.8542e-12 [constexpr]
```

Permittivity vacuum.

2.1.3.3 epsilon_r_1

```
constexpr double epsilon_r_1 = 2.0 [constexpr]
```

Permittivity oil.

2.1.3.4 epsilon_r_2

```
constexpr double epsilon_r_2 = 4.0 [constexpr]
```

Permittivity paper.

2.1.3.5 extra_refinement

```
bool extra_refinement = true
```

true for all cases except Test 1

2.1.3.6 maxlevel

```
constexpr int maxlevel = 6 [constexpr]
```

Level for local refinement.

2.1.3.7 minlevel

```
constexpr int minlevel = 4 [constexpr]
```

Level for global coarsening.

2.1.3.8 N_rhos

```
constexpr size_t N_rhos = 6 [constexpr]
```

Number of points to select for output.

2.1.3.9 NUM_REFINEMENTS

```
constexpr int NUM_REFINEMENTS = 5 [constexpr]
```

Level for global refinement.

2.1.3.10 points

```
std::vector<std::vector<double>> > points {{5e-4,5e-4,0.0},{5e-4,5e-4,z_paper},{5e-4,5e-4,z_oil+z_paper},{5e-
```

Coordinates of the selected points.

2.1.3.11 rho_idx

```
std::vector<size_t> rho_idx
```

2.1.3.12 save_sol

```
constexpr bool save_sol = true [constexpr]
```

If set to 'true' saves data for Paraview visualization.

2.1.3.13 sigma_

```
constexpr double sigma_ = 3.21e-14 [constexpr]
```

Conductivity.

2.1.3.14 T

```
constexpr double T = 5000 [constexpr]
```

Final time of simulation.

2.1.3.15 tau

```
constexpr double tau = 50.0 [constexpr]
```

Time constant for boundary conditions.

2.1.3.16 tol

```
constexpr double tol = 1e-5 [constexpr]
```

Tolerance value for refinement and point selection.

2.1.3.17 tols

```
std::vector<std::vector<double> > tols {{1e-4,1e-4,tol},{1e-4,1e-4,tol},{1e-4,1e-4,tol},{1e-4,1e-4,tol},{1e-
```

Tolerance around the selected points.

2.1.3.18 z_oil

```
constexpr double z_oil = 5e-5 [constexpr]
```

Thickness oil layer.

2.1.3.19 z_paper

```
constexpr double z_paper = 3e-4 [constexpr]
```

Thickness paper layer and butt gaps side length.

2.2 test_3.h

[Go to the documentation of this file.](#)

```
1
2 #include <tmesh_3d.h>
3 #include <simple_connectivity_3d_thin.h> // Loading library with domain geometry
4
5 constexpr int NUM_REFINEMENTS = 5;
6 constexpr int maxlevel = 6;
7 constexpr int minlevel = 4;
8 constexpr double DELTAT = 50.0;
9 constexpr double T = 5000;
10 constexpr double tau = 50.0;
11 constexpr bool save_sol = true;
12
13 // Problem parameters
14 constexpr double epsilon_0 = 8.8542e-12;
15 constexpr double epsilon_r_1 = 2.0;
16 constexpr double epsilon_r_2 = 4.0;
17 constexpr double sigma_ = 3.21e-14;
18 constexpr double z_oil = 5e-5;
19 constexpr double z_paper = 3e-4;
20 constexpr double tol = 1e-5;
21 constexpr size_t N_rhos = 6;
22 std::vector<size_t> rho_idx;
23 std::vector<std::vector<double>>
24   points({{5e-4, 5e-4, 0.0}, {5e-4, 5e-4, z_paper}, {5e-4, 5e-4, z_oil+z_paper}, {5e-4, 5e-4, z_oil+2*z_paper}, {5e-4, 5e-4, 2*z_oil+2*z_paper}});
25 std::vector<std::vector<double>>
26   tols({{1e-4, 1e-4, tol}, {1e-4, 1e-4, tol}, {1e-4, 1e-4, tol}, {1e-4, 1e-4, tol}, {1e-4, 1e-4, tol}, {1e-4, 1e-4, tol}});
27 bool extra_refinement = true;
28 static int
29 uniform_refinement (tmesh_3d::quadrant_iterator q)
30 { return NUM_REFINEMENTS; }
31 static int
32 refinement (tmesh_3d::quadrant_iterator quadrant)
33 {
34   int currentlevel = static_cast<int> (quadrant->the_quadrant->level);
35   double zcoord;
36   int retval = 0;
37   for (int ii = 0; ii < 8; ++ii)
38   {
39     zcoord = quadrant->p(2, ii);
40
41     if (zcoord > z_paper - tol || zcoord < 2*z_paper+2*z_oil+tol)
42     {
43       retval = maxlevel - currentlevel;
44       break;
45     }
46   }
47
48   if (currentlevel >= maxlevel)
49     retval = 0;
50
51   return retval;
52 }
53 static int
54 coarsening (tmesh_3d::quadrant_iterator quadrant)
55 {
56   int currentlevel = static_cast<int> (quadrant->the_quadrant->level);
```

```

64 double xcoord,ycoord,zcoord;
65 int retval = currentlevel - minlevel;
66 for (int ii = 0; ii < 8; ++ii)
67 {
68     xcoord = quadrant->p(0, ii);
69     ycoord = quadrant->p(1, ii);
70     zcoord = quadrant->p(2, ii);
71
72     if (fabs(zcoord - z_paper) < tol || fabs(zcoord - z_paper-z_oil) < tol || fabs(zcoord -
2*z_paper-z_oil) < tol || fabs(zcoord - 2*z_paper-2*z_oil) < tol ||
73         (xcoord<z_paper && fabs(ycoord-5e-4+z_paper/2)<tol && zcoord>5e-4-z_paper/2 &&
zcoord<5e-4+z_paper/2) ||
74         (xcoord<z_paper && fabs(ycoord-5e-4-z_paper/2)<tol && zcoord>5e-4-z_paper/2 &&
zcoord<5e-4+z_paper/2) ||
75         (fabs(xcoord-z_paper)<tol && ycoord>5e-4-z_paper/2 && ycoord<5e-4+z_paper/2 &&
zcoord>5e-4-z_paper/2 && zcoord<5e-4+z_paper/2))
76     {
77         retval = 0;
78         break;
79     }
80 }
81
82 if (currentlevel <= minlevel)
83     retval = 0;
84
85 return (retval);
86 }
87
88 double epsilon_fun(const double & x, const double & y, const double & z)
89 {
90     if ((z > z_paper && z<z_paper+z_oil) || (z>2*z_paper+z_oil && z<2*(z_paper+z_oil)))
91         return epsilon_0 * epsilon_r_1;
92
93     if ((z > z_paper+z_oil && z<2*z_paper+z_oil) && x<z_paper && (y>5e-4-z_paper/2 && y<5e-4+z_paper/2))
94         return epsilon_0 * epsilon_r_1;
95
96     return epsilon_0 * epsilon_r_2;
97 }
98
99 double sigma_fun(const double & x, const double & y, const double & z)
100 {return sigma_ * DELTAT;}
101
102 std::vector<size_t> find_idx(tmsh_3d &tmsh,std::vector<std::vector<double>>
&points,std::vector<std::vector<double>> &tols, const size_t &N_rhos)
103 {
104     std::vector<size_t> id(N_rhos,0);
105
106     for (size_t i = 0; i < N_rhos; i++) {
107         bool found = false;
108
109         for (auto quadrant = tmsh.begin_quadrant_sweep ();
110             quadrant != tmsh.end_quadrant_sweep ();
111             ++quadrant){
112
113             for (int ii = 0; ii < 8; ++ii) {
114
115                 if (fabs(quadrant->p(0,ii)-points[i][0])<tols[i][0] &&
116                     fabs(quadrant->p(1,ii)-points[i][1])<tols[i][1] && fabs(quadrant->p(2,ii)-points[i][2])<tols[i][2]) {
117                     id[i] = quadrant->t(ii);
118                     found = true;
119                     std::cout << "Point " << i+1 << ": x= " << quadrant->p(0,ii) << ", y= " << quadrant->p(1,ii) << ",
120                     z= " << quadrant->p(2,ii) << std::endl;
121                     break;
122                 }
123             }
124
125             if (found)
126                 break;
127         }
128         if(!found)
129             std::cout << "Node " << i+1 << " not found in current rank" << std::endl;
130     }
131     return id;
132 }

```

2.3 HVDC_main.cpp File Reference

```

#include <iostream>
#include <fstream>
#include <cmath>
#include <algorithm>
#include <octave_file_io.h>

```

```
#include <bim_distributed_vector.h>
#include <bim_sparse_distributed.h>
#include <bim_timing.h>
#include <mumps_class.h>
#include <quad_operators_3d.h>
#include <test_4.h>
```

Functions

- int `main` (int argc, char **argv)

2.3.1 Detailed Description

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Version

0.1

2.3.2 LICENSE

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2.3.3 Function Documentation

2.3.3.1 main()

```
int main (
    int argc,
    char ** argv )
{
    using q1_vec = q1_vec<distributed_vector>;
    /*
    45 Manegement of solutions ordering: ord0-> phi          ord1->rho
    46 Equation ordering: ord0->diffusion-reaction equation  ord1->continuity equation
    47 */
    48 ordering
    49   ord0 = [] (tmesh_3d::idx_t gt) -> size_t { return dof_ordering<2, 0> (gt); },
    50   ord1 = [] (tmesh_3d::idx_t gt) -> size_t { return dof_ordering<2, 1> (gt); };
    51
    52 // Initialize MPI
```

```

53 MPI_Init (&argc, &argv);
54 int rank, size;
55 MPI_Comm_rank (MPI_COMM_WORLD, &rank);
56 MPI_Comm_size (MPI_COMM_WORLD, &size);
57
58 // Generate the mesh in 3d
59 tmsh_3d tmsh;
60 tmsh.read_connectivity (simple_conn_p, simple_conn_num_vertices,
61                         simple_conn_t, simple_conn_num_trees);
62
63 //Uniform refinement
64 int recursive = 1;
65 tmsh.set_refine_marker (uniform_refinement);
66 tmsh.refine (recursive);
67
68 //In test 1 we only have uniform refinement, in all other cases we perform additional refinement
69 if (extra_refinement)
70 {
71     tmsh.set_refine_marker(refinement);
72     tmsh.refine (recursive);
73
74     tmsh.set_coarsen_marker(coarsening);
75     tmsh.coarsen(recursive);
76 }
77
78 tmsh_3d::idx_t gn_nodes = tmsh.num_global_nodes ();
79 tmsh_3d::idx_t ln_nodes = tmsh.num_owned_nodes ();
80 tmsh_3d::idx_t ln_elements = tmsh.num_local_quadrants ();
81
82 // Allocate linear solver
83 mumps *lin_solver = new mumps ();
84
85 // Allocate initial data container
86 ql_vec sold (ln_nodes * 2);
87 sold.get_owned_data ().assign (sold.get_owned_data ().size (), 0.0);
88
89 ql_vec sol (ln_nodes * 2);
90 sol.get_owned_data ().assign (sol.get_owned_data ().size (), 0.0);
91
92 std::vector<double> xa;
93 std::vector<int> ir, jc;
94
95 // Declare system matrix
96 distributed_sparse_matrix A;
97 A.set_ranges (ln_nodes * 2);
98
99 // Buffer for export filename
100 char filename[255]="";
101
102 //Output rho vector
103 size_t N_timesteps = (size_t) (ceil(T/DELTAT)+1);
104 std::vector<std::vector<double>> rho_out(N_timesteps, std::vector<double>(N_rhos+1));
105
106 // Compute coefficients
107
108 // diffusion
109 std::vector<double> epsilon (ln_elements, 0.);
110 std::vector<double> sigma (ln_elements, 0.);
111 ql_vec zero (ln_nodes);
112
113 // reaction
114 std::vector<double> delta0 (ln_elements, 0.);
115 std::vector<double> delta1 (ln_elements, 0.);
116 ql_vec zeta0 (ln_nodes);
117 ql_vec zeta1 (ln_nodes);
118
119 // rhs
120 std::vector<double> f0 (ln_elements, 0.);
121 std::vector<double> f1 (ln_elements, 0.);
122 ql_vec g0 (ln_nodes);
123 ql_vec g1 (ln_nodes);
124
125 // Initialize constant (in time) parameters and initial data
126 for (auto quadrant = tmsh.begin_quadrant_sweep ();
127      quadrant != tmsh.end_quadrant_sweep ();
128      ++quadrant)
129 {
130     double xx{quadrant->centroid(0)}, yy{quadrant->centroid(1)}, zz{quadrant->centroid(2)};
131
132     epsilon[quadrant->get_forest_quad_idx ()] = epsilon_fun(xx,yy,zz);
133     sigma[quadrant->get_forest_quad_idx ()] = sigma_fun(xx,yy,zz);
134
135     delta0[quadrant->get_forest_quad_idx ()] = -1.0;
136     delta1[quadrant->get_forest_quad_idx ()] = 1.0;
137     f0[quadrant->get_forest_quad_idx ()] = 0.0;
138     f1[quadrant->get_forest_quad_idx ()] = 1.0;
139 }

```



```

140     for (int ii = 0; ii < 8; ++ii)
141     {
142         if (! quadrant->is_hanging (ii))
143         {
144             zero[quadrant->gt (ii)] = 0.;
145             zeta0[quadrant->gt (ii)] = 1.0;
146             zeta1[quadrant->gt (ii)] = 1.0;
147             g0[quadrant->gt (ii)] = 0.;
148
149             double zz=quadrant->p(2,ii);
150             sold[ord0(quadrant->gt (ii))] = 0.0;
151             sold[ord1(quadrant->gt (ii))] = 0.0;
152             sol[ord0(quadrant->gt (ii))] = 0.0;
153             sol[ord1(quadrant->gt (ii))] = 0.0;
154         }
155         else
156             for (int jj = 0; jj < quadrant->num_parents (ii); ++jj)
157             {
158                 zero[quadrant->gparent (jj, ii)] += 0.;
159                 zeta0[quadrant->gparent (jj, ii)] += 0.;
160                 zeta1[quadrant->gparent (jj, ii)] += 0.;
161                 g0[quadrant->gparent (jj, ii)] += 0.;
162
163                 sold[ord0(quadrant->gparent (jj, ii))] += 0.;
164                 sold[ord1(quadrant->gparent (jj, ii))] += 0.;
165             }
166     }
167 }
168
169 bim3a_solution_with_ghosts (tmsh, sold, replace_op, ord0, false);
170 bim3a_solution_with_ghosts (tmsh, sold, replace_op, ord1);
171
172 zero.assemble (replace_op);
173 zeta0.assemble (replace_op);
174 zeta1.assemble (replace_op);
175 g0.assemble (replace_op);
176
177 // Save initial conditions
178 sprintf(filename, "model_0_u_0000");
179 tmsh.octbin_export (filename, sold, ord0);
180 sprintf(filename, "model_0_v_0000");
181 tmsh.octbin_export (filename, sold, ord1);
182
183 int count = 0;
184
185 //Choosing the indices for the nodes corresponding to the vales of rho of interest
186 if (rank == 0) {
187     rho_out[0]=std::vector<double>(N_rhos+1,0.0);
188     rho_idx = find_idx(tmsh,points,tols,N_rhos);
189 }
190
191 // Time cycle
192 for( double time = DELTAT; time <= T; time += DELTAT)
193 {
194     count++;
195
196     // Define boundary conditions
197     dirichlet_bcs3 bcs0, bcs1;
198     bcs0.push_back (std::make_tuple (0, 4, [(double x, double y, double z){return 0.0;}])); //bottom
199     bcs0.push_back (std::make_tuple (0, 5, [time](double x, double y, double z){return 1.5e4 * (1 -
exp(-time/tau));})); //top
200
201     // Print current time
202     if(rank==0)
203         std::cout<<"TIME= "<<time<<std::endl;
204
205     // Reset containers
206     A.reset ();
207     sol.get_owned_data ().assign (sol.get_owned_data ().size (), 0.0);
208     sol.assemble (replace_op);
209
210     // Initialize non constant (in time) parameters
211     for (auto quadrant = tmsh.begin_quadrant_sweep ();
212          quadrant != tmsh.end_quadrant_sweep ();
213          ++quadrant)
214     {
215         for (int ii = 0; ii < 8; ++ii)
216             if (! quadrant->is_hanging (ii))
217                 g1[quadrant->gt (ii)] = sold[ord1(quadrant->gt (ii))];
218
219         else
220             for (int jj = 0; jj < quadrant->num_parents (ii); ++jj)
221                 g1[quadrant->gparent (jj, ii)] += 0.;
222     }
223
224     g1.assemble(replace_op);
225

```

```

226 // advection_diffusion
227 bim3a_advection_diffusion (tmsh, epsilon, zero, A, true, ord0, ord0);
228 bim3a_advection_diffusion (tmsh, sigma, zero, A, true, ord1, ord0);
229
230 // reaction
231 bim3a_reaction (tmsh, delta0, zeta0, A, ord0, ord1);
232 bim3a_reaction (tmsh, delta1, zeta1, A, ord1, ord1);
233
234 //rhs
235 bim3a_rhs (tmsh, f0, g0, sol, ord0);
236 bim3a_rhs (tmsh, f1, g1, sol, ord1);
237
238 //boundary conditions
239 bim3a_dirichlet_bc (tmsh, bcs0, A, sol, ord1, ord0, false);
240
241 // Communicate matrix and RHS
242 A.assemble ();
243 sol.assemble ();
244
245 // Solver analysis
246 lin_solver->set_lhs_distributed ();
247 A.ajj (xa, ir, jc, lin_solver->get_index_base ());
248 lin_solver->set_distributed_lhs_structure (A.rows (), ir, jc);
249 std::cout << "lin_solver->analyze () return value = " << lin_solver->analyze () << std::endl;
250
251 // Matrix update
252 A.ajj_update (xa, ir, jc, lin_solver->get_index_base ());
253 lin_solver->set_distributed_lhs_data (xa);
254
255 // Factorization
256 std::cout << "lin_solver->factorize () = " << lin_solver->factorize () << std::endl;
257
258 // Set RHS data
259 lin_solver->set_rhs_distributed (sol);
260
261 // Solution
262 std::cout << "lin_solver->solve () = " << lin_solver->solve () << std::endl;
263
264 // Copy solution
265 ql_vec result = lin_solver->get_distributed_solution ();
266 for (int idx = sold.get_range_start (); idx < sold.get_range_end (); ++idx)
267     sold (idx) = result (idx);
268 sold.assemble (replace_op);
269
270 // Save solution
271 if (save_sol == true)
272 {
273     sprintf(filename, "model_0_u_%.4d", count);
274     tmsh.octbin_export (filename, sold, ord0);
275     sprintf(filename, "model_0_v_%.4d", count);
276     tmsh.octbin_export (filename, sold, ord1);
277 }
278
279 // Save rho values
280 if (rank == 0)
281 {
282     std::vector<double> temp(N_rhos+1);
283     temp[0] = time;
284     for (size_t i=1; i < N_rhos+1; i++)
285         temp[i] = sold[ord1(rho_idx[i-1])];
286
287     rho_out[count] = temp;
288 }
289 }
290
291 // Print file with rho values
292 if (rank == 0)
293 {
294     std::ofstream outFile("Arho.txt");
295     outFile << N_rhos+1 << "\t" << N_timesteps << "\n";
296
297     for (const auto &e : rho_out) {
298         for (size_t i=0; i < N_rhos+1; i++)
299             outFile << e[i] << "\t";
300         outFile << "\n";
301     }
302 }
303
304 // Close MPI and print report
305 MPI_Barrier (MPI_COMM_WORLD);
306
307 // Clean linear solver
308 lin_solver->cleanup ();
309
310 MPI_Finalize ();
311
312 return 0;

```

```
313 }
```


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