source/slae.hpp

```
1
   #pragma once
3
   #include "utils.hpp"
4
   #include <cassert>
 5
6
7
8
   // Backwards gaussian elimination. O(N^2) complexity.
9
10
   // Assumes 'R' to be upper-triangular matrix.
11
   //
12
   inline Vector backwards_gaussian_elimination(const Matrix& R, Vector rhs) {
13
        assert(R.rows() == R.cols());
14
        assert(R.rows() == rhs.rows());
15
        for (Idx i = R.rows() - 1; i >= 0; --i) {
16
17
            for (Idx j = i + 1; j < R.cols(); ++j) rhs(i) -= R(i, j) * rhs(j);
18
            rhs(i) /= R(i, i);
19
        }
20
21
        return rhs;
22
   }
23
24
25
   // Forward gaussian elimination. O(N^3) complexity.
26
   //
27
   // Partial pivoting.
28
29
   inline void partial_piv_forward_gaussian_elimination(Matrix& A, Vector& rhs) {
30
        assert(A.rows() == A.cols());
31
        assert(A.rows() == rhs.rows());
32
33
        for (Idx i = 0; i < A.rows(); ++i) {
34
            // Partial pivot
35
            Idx i max = i;
            for (Idx ii = i; ii < A.rows(); ++ii)</pre>
36
37
                if (std::abs(A(ii, i)) > std::abs(A(i_max, i))) i_max = ii;
38
39
            if (i != i max) {
40
                // Swap rows (matrix)
                const Vector tmp A
41
                                     = A.row(i);
42
                A.row(i)
                                      = A.row(i max);
43
                A.row(i max)
                                      = tmp A;
44
                // Swap rows (rhs)
45
                const double tmp_rhs = rhs(i);
46
                rhs(i)
                                     = rhs(i max);
47
                rhs(i_max)
                                      = tmp_rhs;
48
            }
49
50
            // Elimination (normalize current row)
51
            const double factor = 1. / A(i, i);
52
            for (Idx j = i; j < A.cols(); ++j) A(i, j) *= factor;
53
            rhs(i) *= factor;
54
55
            // Elimination (substract current row from all the rows below)
56
            for (Idx k = i + 1; k < A.rows(); ++k) {
```

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```
57
                const double first = A(k, i);
58
                for (Idx j = i; j < A.cols(); ++j) A(k, j) -= first * A(i, j);
59
                rhs(k) -= first * rhs(i);
60
            }
61
       }
62
63
   // Forward + backwards gaussian elimination. O(N^3) complexity.
64
65
   // Partial pivoting. Jacobi preconditioner.
66
67
   //
   inline Vector partial piv gaussian elimination(Matrix A, Vector rhs) {
68
       // Jacobi preconditioner
69
70
       //
       // When testing N = 4 one of the (A - lambda I) matrices with epsilon = 0.1
71
   was almost degenerate (det = 1e-7),
       // without preconditioning it went into 'nan's, Jacobi was just good enough to
72
   prevent this.
73
       //
       // N = 10 case is luckier and doesn't require this, but why not do it
74
   regardless.
75
       //
76
       Matrix preconditioner
                                  = Matrix::Zero(A.rows(), A.cols());
77
       preconditioner.diagonal() = A.diagonal();
78
       Α
                                  = preconditioner * A;
79
                                  = preconditioner * rhs;
       rhs
80
81
       // Gaussian elimination
82
       partial_piv_forward_gaussian_elimination(A, rhs);
83
        return backwards gaussian elimination(A, rhs);
84 }
```

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