

## Schedule

Time	Descrption	Туре
9.00 - 9.05	Introduciton	Lecture
9.05 - 9.15	Setting up development environment	Lab/Interactive
9.15 - 10.00	Working with matrices and vectors	Lecture
10.00 - 10.20	Coffee break	
10.20 - 11.00	Working with exercises	Lab
11.00 - 12.00	Advanced matrix operations	Lecture
12.00 - 13.00	Lunch	
13.00 - 13.40	Working with exercises	Lab
13.40 - 14.00	Best practice and integration	Lecture
14.00 - 14.30	Working with exercises	Lab
14.30 - 14.45	Coffee break	Lecture
14.45 - 15.15	Using Eigen in Parallel applications	Lecture

#### Course page

- Main page
  - https://array-computing-with-eigen.readthedocs.io/en/latest/index.html
- Exercises
  - https://array-computing-witheigen.readthedocs.io/en/latest/exercises.html

#### Array Computing with Eigen Contents:

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Setup

Schedule

Links

Exercises

#### Array Computing with Eigen





This is a course on array computing with the C++ library Eigen. The course is intended for students with a basic knowledge of C++ programming and linear algebra.

Jonas Lindemann, 2025





## Why Not Build Your Own Matrix Library

- Arrays are crucial for scientific computing applications
- In C++, it might be tempting to create your own matrix library
- However, this is often not recommended because:
  - Existing libraries are well-tested
  - Existing libraries are highly optimized
  - Creating your own introduces unnecessary complexity



## What is Eigen

- One of the most popular libraries for linear algebra in C++
- A header-only library (no linking required)
- Fast and easy to use
- Supports optimized packages like BLAS and LAPACK
- Can significantly speed up computational tasks
  - Some operations have support for OpenMP
  - Using optimized libraries will provide additional speed.



## Setting up Eigen in Your Project

- Eigen is header-only, so no linking needed
  - Can still need linking for optimized BLAS / LAPACK libraries
- Simply include the relevant header:

#include <Eigen/Dense>

- The Dense module includes all basic matrix operations
- Most commonly used module for linear algebra tasks
- Other specialized modules are available for specific needs

## Getting Started with Eigen

- Easy to integrate into existing C++ projects
- No external dependencies required
- Compatible with modern C++ standards
- Excellent documentation available
- Active community support

## Setting Up Your Eigen Development Environment

## **Environment Setup**

- COSMOS at LUNARC (HPC environment)
- Windows
- Linux (Ubuntu)
- macOS

## Using COSMOS at LUNARC

- Remote desktop environment: LUNARC Documentation
- SSH access: <u>SSH Login Documentation</u>

## LUNARC: Loading the Environment

#### Load required modules:

```
module load foss/2024a
module load Eigen
module load Cmake
```

#### Verify installation:

```
g++ --version # Should show GCC 13.3.0
cmake --version # Should show CMake 3.29.3
```

## LUNARC: Testing Eigen

Create a test file ex0.cpp:

```
#include <iostream>
#include <Eigen/Dense>

int main()
{
    Eigen::Matrix3d m = Eigen::Matrix3d::Random();
    std::cout << "Here is the matrix m:" << std::endl;
    std::cout << m << std::endl;
    return 0;
}</pre>
```

## LUNARC: Compiling and Running

```
Compile:
 g++ ex0.cpp -o ex0
Run:
 ./ex0
Expected output:
 Here is the matrix m:
 0.680375 0.59688 - 0.329554
 -0.211234 0.823295 0.536459
 0.566198 -0.604897 -0.444451
```

## Ubuntu Linux Setup

#### Install required packages:

```
sudo apt-get update
sudo apt-get install g++
sudo apt-get install cmake
sudo apt-get install libeigen3-dev
```

#### Verify installation:

```
g++ --version cmake --version
```

## Ubuntu Linux: Testing Eigen

Create the same test file ex0. cpp as before. Compile with include path:

```
g++ ex0.cpp -I/usr/include/eigen3 -o ex0
```

Run:

./ex0

The output should match the previous example.

#### macOS Setup

Install required packages using Homebrew:

```
brew install gcc
brew install cmake
brew install eigen
```

## macOS: Testing Eigen

Create the same test file ex0. cpp as before.

Compile with include path:

```
g++ -std=c++11 -I/usr/local/include/eigen3 ex0.cpp -o ex0
```

Run:

./ex0

The output should match the previous examples.

## Troubleshooting Tips

- Check compiler version compatibility
- Verify include paths are correct
- Ensure Eigen headers are properly installed
- For library issues, confirm environment variables are set correctly

# 



## Matrix and Vector Types

- All Eigen classes are template classes (work with different data types)
- Most common data types: float, double, and int
- All classes defined in the **Eigen** namespace
- Best practice: use Eigen:: prefix instead of
  - using namespace Eigen
- In many of my example I will use without the prefix to make the code more readable.

#### Matrix Declaration

#### **Generic form:**

```
// 3x3 matrix of doubles
Eigen::Matrix<double, 3, 3> A;
```

#### Using convenient typedefs:

```
Eigen::Matrix3d B; // Same as above
```

#### Vector Declaration

#### **Generic form:**

```
// 3x1 vector of doubles
Eigen::Vector<double, 3> v;
```

#### Using convenient typedefs:

```
Eigen::Vector3d w; // Same as above
```

Note: Vectors are matrices with one dimension fixed to 1

#### Initialization

- Newly declared matrices contain random values
- Methods to initialize:
  - A.setZero() Initialize to zeros
  - A.setOnes() Initialize to ones
  - << operator Set specific values</li>

### Vectors - Special Properties

Initialize with constructor:

```
Eigen:: Vector3d v(1, 2, 3);
```

Initialize with the << operator:</li>

```
Eigen::Vector3d w;
w << 1, 2, 3;</pre>
```

Special initialization methods:

```
v.setLinSpaced(3, 1, 2); // Creates: [1, 1.5, 2]
w.setRandom(); // Random values
```

#### Column and Row Access

```
// Column insertion
D.col(0) << 1, 2, 3;
D.col(1) << 4, 5, 6;
D.col(2) << 7, 8, 9;
// Row insertion
D.row(0) << 1, 2, 3;
D.row(1) << 4, 5, 6;
D.row(2) << 7, 8, 9;
```

## Common Typedefs

#### Fixed-size matrices:

```
Matrix2d, Matrix3d, Matrix4d // double
Matrix2f, Matrix3f, Matrix4f // float
Matrix2i, Matrix3i, Matrix4i // int
```

#### Dynamic-size matrices:

```
MatrixXd // double
MatrixXf // float
MatrixXi // int
```

## Common Vector Typedefs

#### Fixed-size vectors:

```
Vector2d, Vector3d, Vector4d // double Vector2f, Vector3f, Vector4f // float Vector2i, Vector3i, Vector4i // int
```

#### Dynamic-size vectors:

```
VectorXd // double
VectorXf // float
VectorXi // int
```

# 

## Fixed vs Dynamic Size

```
Fixed size (known at compile time):
 // 3x3 matrix, size fixed at compile time
 Eigen::Matrix3d A;
Dynamic size (determined at runtime):
 Eigen::Matrix<double, Eigen::Dynamic, Eigen::Dynamic> F;
 F.resize(3, 3); // Set dimensions at runtime
 // Shorthand
 Eigen::MatrixXd G(3, 3);
```

## Resizing Matrices

- resize(rows, cols) changes matrix dimensions
- If the total number of elements stays the same, data is preserved
- If the total elements change, data is lost and must be reinitialized

```
A_dyn_resize(1, 9); // Preserves data if A_dyn was 3x3
A_dyn_resize(6, 6); // Changes total elements, data is lost
```

#### Row Vectors

```
// 1x3 vector
Eigen::RowVector3d r(1.0, 2.0, 3.0);
// Output: 1 2 3
// 3x1 vector
Eigen::Vector3d s(1.0, 2.0, 3.0);
// Output:
// 1
// 2
// 3
```

## Matrix Operations

- Addition: A + B
- Scalar multiplication: A \* 3.0
- Matrix multiplication: A \* B
- Add scalar to all elements: E + Matrix3d::Constant(1.0)
- Element-wise operations: E\_array() + 3.0

#### Matrix Transformations

- Transpose: A.transpose()
- Inverse: A.inverse()
- Component-wise operations:

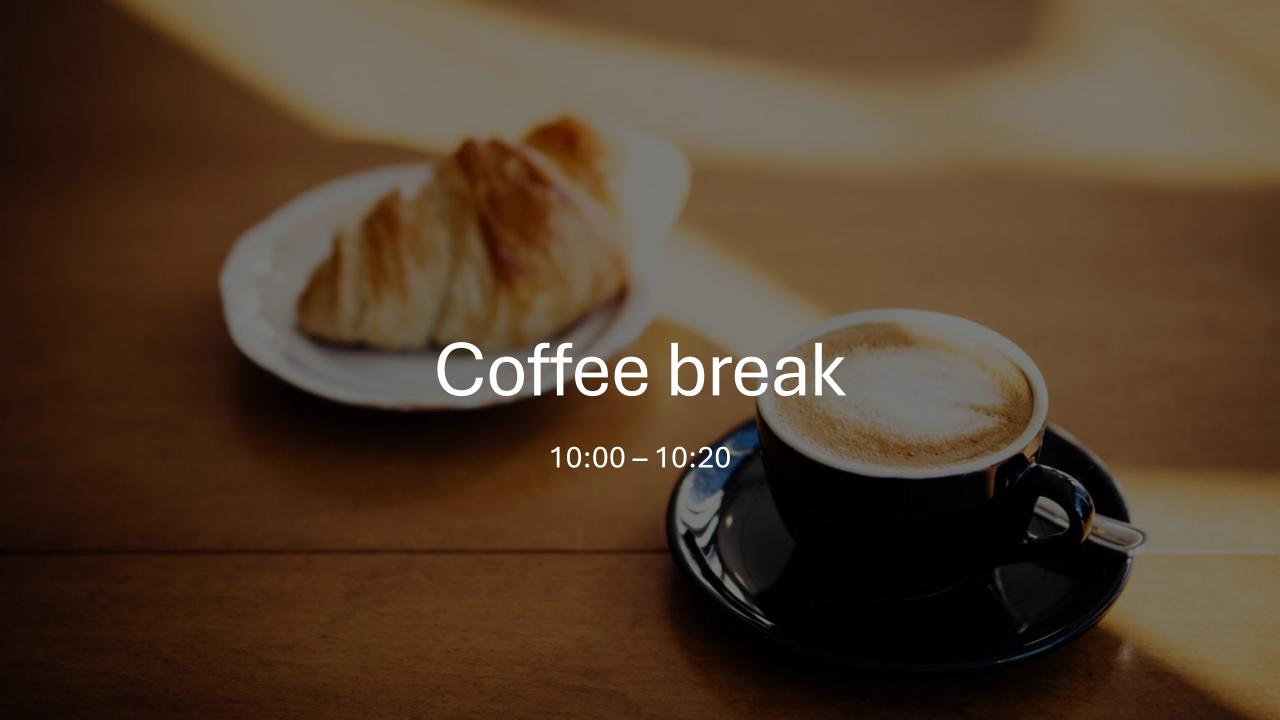
```
Vector3d x(1, 4, 9);
auto y = x.cwiseSqrt(); // [1, 2, 3]
// Or
auto w = z.array().sqrt();
```

# Vector Operations

- Dot product: s.dot(t)
- Cross product: s.cross(t)
- Component-wise operations: x . cwiseSqrt()

# Reduction Operations

- K.sum() Sum of all elements
- K.prod() Product of all elements
- K.mean() Mean of all elements
- K.norm() Euclidean norm
- K.maxCoeff() Maximum value
- K.minCoeff() Minimum value
- K.trace() Trace of the matrix
- K.diagonal() Diagonal elements
- K.determinant() Determinant







# Reshaping Matrices

- reshaped (rows, cols) method allows changing matrix dimensions
- Returns a view into the original matrix (not a copy)
- Changes to the reshaped matrix affect the original Matrix3d A;

```
A << 1, 2, 3,
4, 5, 6,
7, 8, 9;
```

```
auto B = A.reshaped(1, 9); // Result: 1 2 3 4 5 6 7 8 9
```

### Reshaping Considerations

- Data is stored in column-major order in Eigen
- Self-assignment requires \_eval() to force evaluation:
   C = C\_reshaped(1, 9)\_eval();
- Can combine with \_transpose():
   MatrixXd D = C\_reshaped(1, 9)\_transpose();

# Slicing and Indexing

Row and column access:

```
// Set values in row 3
A.row(3) << 1, 2, 3, 4, 5, 6, 7, 8, 9, 10;
// Set values in column 3
A.col(3) << 1, 2, 3, 4, 5, 6, 7, 8, 9, 10;
// Set all elements in column 1 to 1
A.col(1).setOnes();</pre>
```

# Range Selection with seq()

Select ranges of rows and columns:

```
// Select block from rows 3-5, columns 3-5
B(seq(3, 5), seq(3, 5)).setConstant(1);

// Select every other row/column from 0-9
B(seq(0, 9, 2), seq(0, 9, 2)).setConstant(2);
```

### Special Selectors

```
// Set all elements in the last column to 3
B(all, last).setConstant(3);

// Set all elements in the second-to-last
// column to 4
B(all, last - 1).setConstant(4);
```

#### Index-Based Selection

Use std::vector of indices to select specific rows and columns:

```
vector<int> idx = { 1, 3, 4, 6, 7, 9 };

// Select submatrix using index vector
auto D = C(idx, idx);
```

# Linear System Solving - Small Matrices

For small matrices (up to 4x4), using inverse is acceptable:

```
Matrix3d A;
A.setRandom();
Vector3d b;
b.setRandom();

// Solve Ax = b
Vector3d x = A.inverse() * b;
```

# Linear System Solving - Larger Matrices

For larger matrices, use decomposition methods:

```
MatrixXd A(10, 10);
A setRandom();
VectorXd b(10);
b_setRandom();
// Solve using QR decomposition
VectorXd x = A.colPivHouseholderQr().solve(b);
// Check error
double error = (A * x - b).norm();
```

# Matrix Decompositions

Different decompositions for different matrix types:

- colPivHouseholderQr() General matrices (robust)
- fullPivLu() General matrices (most stable, slower)
- ldlt() Symmetric matrices
- householderQr() General matrices (fastest, less accurate)

### Reusing Decompositions

Create decomposition objects for reuse with multiple right-hand sides:

```
// Create decomposition once
FullPivLU<MatrixXd> lu(A);

// Solve for multiple right-hand sides
VectorXd x1 = lu.solve(b1);
VectorXd x2 = lu.solve(b2);

// Solve systems with multiple right-hand sides at once
MatrixXd X = lu.solve(B); // B has multiple columns
```





# Returning Matrices from Functions

- Prefer returning Eigen matrices by value
- C++ return value optimization (RVO) prevents unnecessary copying
- Simple and effective approach:

```
MatrixXd foo()
{
    MatrixXd A(10, 10);
    A.setRandom();
    return A; // Efficient due to RVO
}
int main()
{
    MatrixXd B = foo(); // No unnecessary copying
}
```

### Passing Matrices to Functions

- Pass matrices as const references to avoid copying
- Use the const keyword to indicate the matrix won't be modified

```
void bar(const MatrixXd& A) // Pass by reference
{
   cout << A << endl;
}
int main()
{
   MatrixXd B(10, 10);
   B.setRandom();
   bar(B); // No copying occurs
}</pre>
```

# Implementing Functions with Eigen

**General Rule:** - Pass Eigen matrices and vectors by reference - Return Eigen matrices and vectors by value

```
// Example: Function that creates and returns a matrix
MatrixXd hooke(TAnalysisType ptype, double E, double v)
{
    MatrixXd D;
    // ... matrix construction ...
    return D; // Return by value
}
```

### **Example: Function Implementation**

```
Matrix4d bar2e(const Vector2d& ex, const Vector2d& ey, const
Vector2d& ep)
   // Parameters passed by const reference
    double E = ep(0);
   double A = ep(1);
    double L = sqrt(pow(ex(1)-ex(0),2)+pow(ey(1)-ey(0),2));
   // ... calculations ...
    Matrix4d Ke = G_transpose()*Ke_loc*G;
    return Ke; // Return by value
```

# Accessing Raw Data

- Use .data() method to get pointer to raw data
- Useful for interfacing with C-style APIs

```
MatrixXd A(10, 10);
A.setRandom();

// Get pointer to underlying data
double* data = A.data();

// Access elements directly
for (int i = 0; i < A.size(); i++) {
    cout << data[i] << " ";
}</pre>
```

### Creating 2D Array Views

- Eigen stores data as 1D array internally
- Create array of pointers for 2D access:

### Integrating with Other Libraries

• When working with libraries requiring 2D C-style arrays:

```
void foo(double** data, int rows, int cols) {
    // Function expecting 2D C-style array
// Setup for integration
double** data2D = new double*[A.rows()];
for (int i = 0; i < A rows(); i++)</pre>
    data2D[i] = A_row(i)_data();
// Call external function
foo(data2D, A_rows(), A_cols());
// Clean up
delete[] data2D;
```

# Dealing with Const Issues

Use const\_cast when necessary for C-style API integration:

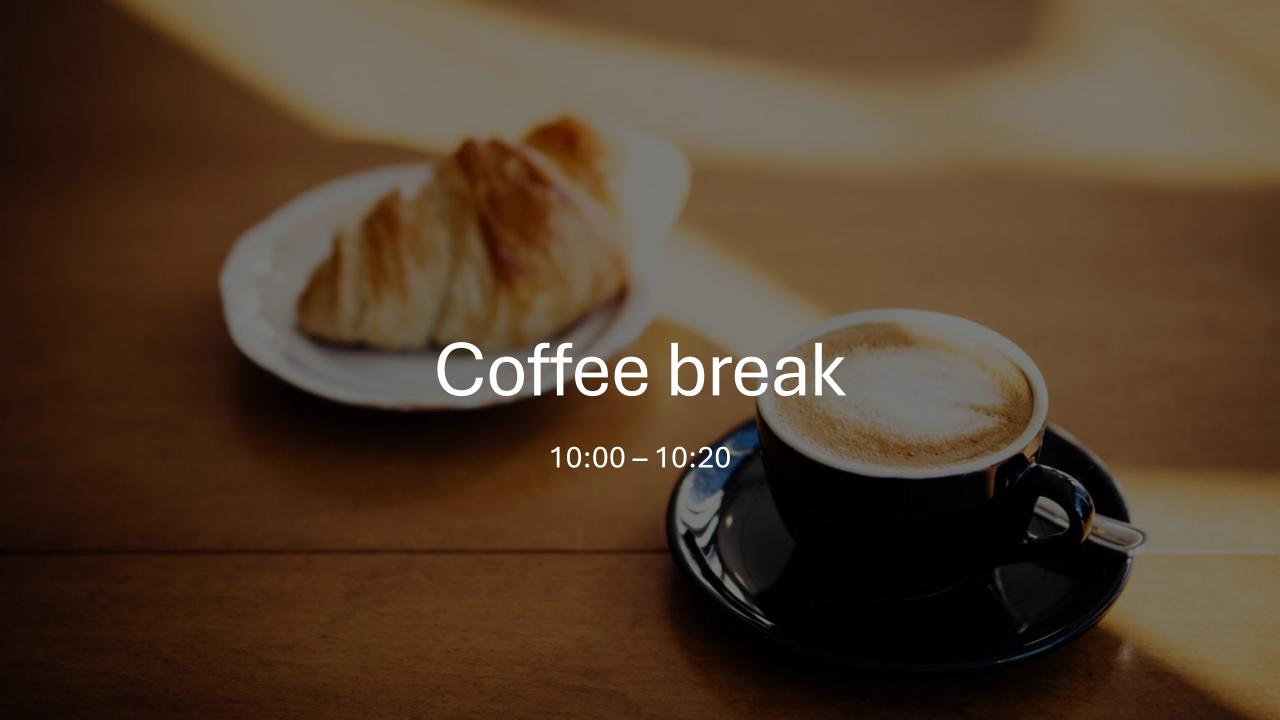
```
// If compiler warns about const correctness
data2D[i] = const_cast<double*>(A.row(i).data());
```

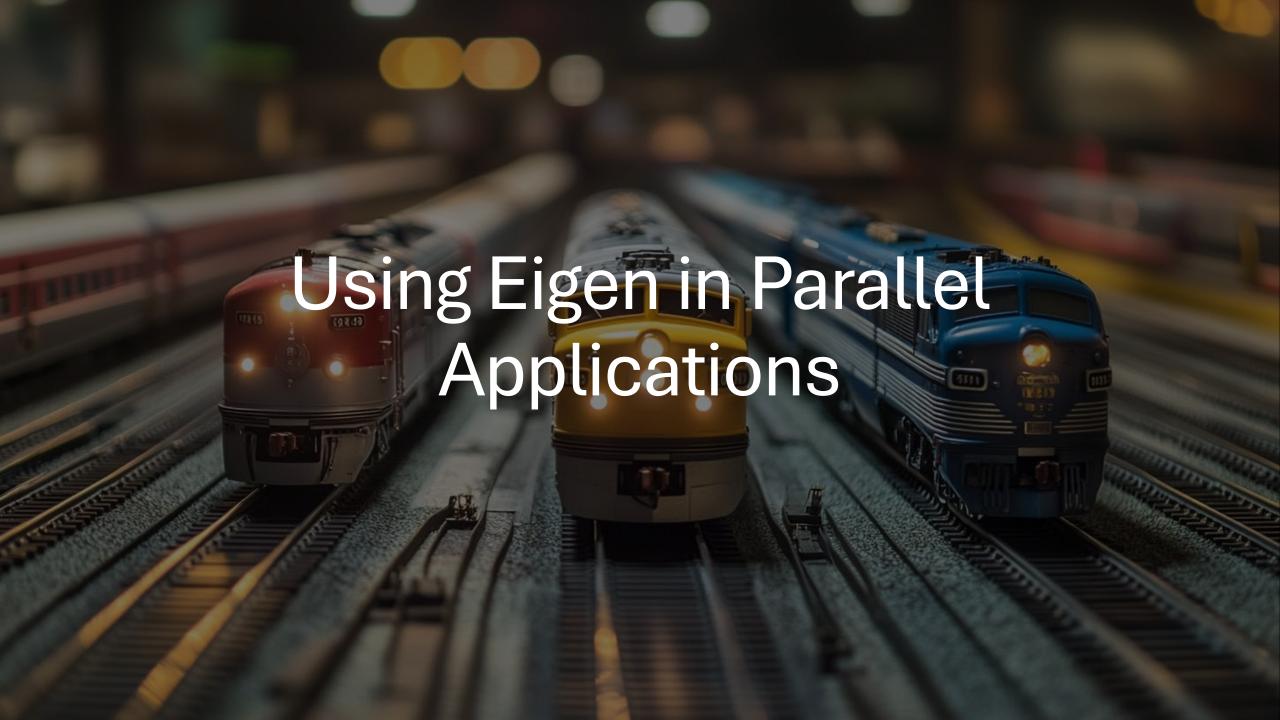
- Only do this when sure the data won't be modified
- Avoid if possible const correctness is a feature, not a bug

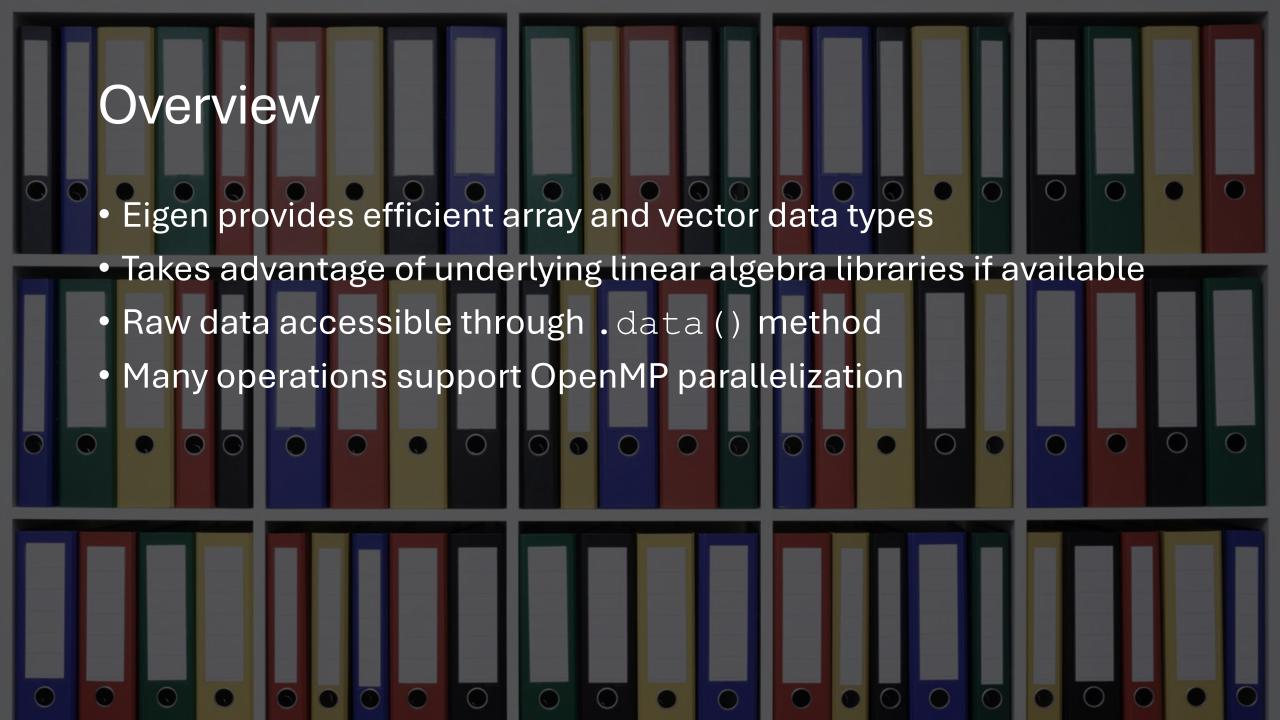
# Summary of Best Practices

- 1. Return matrices by value
- 2. Pass matrices by const reference
- 3. Use data() for raw data access
- 4. Create proper interfaces for C-style libraries
- 5. Maintain const correctness when possible
- 6. Be aware of memory ownership when interfacing









## Eigen and OpenMP Support

Eigen operations that use OpenMP for parallelization:

- general dense matrix matrix products
- PartialPivLU
- row-major-sparse \* dense vector/matrix products
- ConjugateGradient with Lower|Upper as the UpLo template parameter.
- BiCGSTAB with a row-major sparse matrix format.
- LeastSquaresConjugateGradient

## Using Eigen with OpenMP

Simple example: Matrix multiplication using Eigen's built-in OpenMP support

```
// Setting number of threads
omp_set_num_threads(n_threads);

// This operation will automatically use OpenMP
MatrixXd C = A * B;
```

#### OpenMP Performance Scaling

Performance scaling for matrix multiplication (20000×20000):

Threads	Time (seconds)
1	870.111
2	427.386
4	217.594
8	109.546
12	74.313
24	40.574
48	22.657

## Custom Parallelization (Part 1)

Using Eigen for storage with custom OpenMP implementation:

## Custom Parallelization (Part 2)

```
// Get raw pointers to the data
const double* A_data = A.data();
const double* x data = x.data();
double* result data = result.data();
#pragma omp parallel for schedule(static)
for (int i = 0; i < rows; i++) {</pre>
    double sum = 0.0;
    const double* row = A_data + i * cols;
    for (int j = 0; j < cols; j++) {</pre>
        sum += row[j] * x data[j];
    result data[i] = sum;
return result;
```

## Performance Comparison

Custom OpenMP implementation vs. Eigen's built-in (40000×40000):

```
Matrix size: 40000x40000

OpenMP implementation time: 352ms
```

Eigen implementation time: 389ms

Relative error: 7.35186e-15

Custom implementation slightly outperforms Eigen's built-in operation.

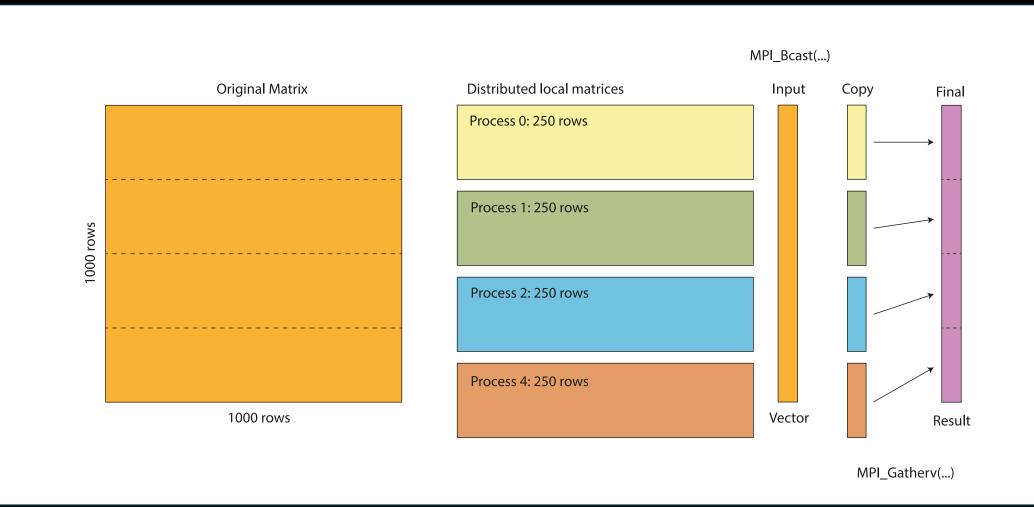
## Using Eigen with MPI

- Eigen is NOT a distributed Matrix library
- MPI for distributed memory parallelism
- Eigen arrays stored in 1D memory layout
- Developer handles data distribution
- Example: matrix-vector multiplication across multiple processes

#### MPI Distribution Strategy

- Matrix divided by rows
- Each process handles a portion of rows
- Vector broadcasted to all processes
- Each process computes partial result

# **MPI Distribution Strategy**



## MPIMatrix Class (Part 1)

```
class MPIMatrix {
private:
    MatrixXd m_localMatrix;
    VectorXd m_localResult;
public:
    MPIMatrix(int r, int c)
        : m_rank(0), m_size(1), m_rows(r), m_cols(c)
        MPI_Comm_rank(MPI_COMM_WORLD, &m_rank);
        MPI_Comm_size(MPI_COMM_WORLD, &m_size);
```

## MPIMatrix Class (Part 2)

```
// Calculate local matrix size
        int localRows = m_rows / m_size;
        if (m_rank < m_rows % m_size) {</pre>
            localRows++;
        m_localMatrix.resize(localRows, m_cols);
        m_localResult.resize(localRows);
    void randomize() {
srand(std::chrono::system_clock::now().time_since_epoch().count());
        m localMatrix.setRandom();
```

## MPIMatrix Class (Part 3)

```
void multiply(const VectorXd& vec) {
   // Local multiplication
   m localResult = m_localMatrix * vec;
void printResult() const {
   // Gather results
   std::vector<int> recvCounts(m_size);
    std::vector<int> displs(m_size);
   // Calculate receive counts and displacements
   int baseCount = m_rows / m_size;
   int remainder = m_rows % m_size;
```

## MPIMatrix Class (Part 4)

```
for (int i = 0; i < m_size; ++i) {</pre>
    recvCounts[i] = baseCount + (i < remainder ? 1 : 0);</pre>
    displs[i] = (i > 0) ? displs[i-1] + recvCounts[i-1] : 0;
// Allocate space for complete result
VectorXd globalResult;
if (m_rank == 0)
    globalResult_resize(m rows);
// Gather all local results to rank 0
MPI_Gatherv(m_localResult.data(), m_localResult.size(), MPI_DOUBLE,
           globalResult.data(), recvCounts.data(), displs.data(),
           MPI_DOUBLE, 0, MPI_COMM_WORLD);
```

## MPIMatrix Class (Part 5)

## MPI Main Function (Part 1)

```
int main(int argc, char** argv) {
    constexpr int MatrixSize = 10000;
    int rank;

MPI_Init(&argc, &argv);
    MPI_Comm_rank(MPI_COMM_WORLD, &rank);

// Create distributed matrix
    MPIMatrix distMatrix(MatrixSize, MatrixSize);
    distMatrix.randomize();
```

## MPI Main Function (Part 2)

```
// Create a random x vector
VectorXd x;
if (rank == 0) {
    std::cout << "Generating random vector x...\n";
    x = VectorXd::Random(MatrixSize);
} else {
    x.resize(MatrixSize);
}
// Broadcast x vector to all processes
MPI_Bcast(x.data(), MatrixSize, MPI_DOUBLE, 0, MPI_COMM_WORLD);</pre>
```

## MPI Main Function (Part 3)

#### MPI Considerations

- For large datasets, avoid gathering all data to one process
- Consider:
  - Writing results to separate files on each rank
  - Using parallel I/O libraries like HDF5
  - Using MPI I/O functions

## Summary: Eigen in Parallel Applications

- Eigen works well with OpenMP and MPI
- Use built-in Eigen parallel operations when possible
- For custom parallelization:
  - Access raw data with .data()
  - Use RowMajor order for row-wise operations
  - Implement your own parallelization with OpenMP/MPI
- Benefits:
  - Simplified memory management
  - Clean, maintainable code
  - Good performance

