

Calcul Scientifique

# Projet de Calcul Scientifique

Élèves :

THEVENET Louis
SABLAYROLLES Guillaume

# 1.

# 1.1.

Matrix dimension	Matrix type	Exec. time for eig (s)	Exec. time for power_v11, (s)
$200 \times 200$	Type 1	9.000e-02	1.510e+00
$400 \times 400$	Type 1	4.000e-02	1.831e+01
$600 \times 600$	Type 1	6.000 e-02	6.021e+01
$200 \times 200$	Type 2	3.000 e-02	3.000e-02
$400 \times 400$	Type 2	4.000e-02	4.000e-02
$600 \times 600$	Type 2	7.000e-02	1.700e-01
$200 \times 200$	Type 3	1.000e-02	5.000e-02
$400 \times 400$	Type 3	3.000e-02	5.200e-01
$600 \times 600$	Type 3	7.000e-02	1.270e+00
$200 \times 200$	Type 4	2.000e-02	1.670e+00
$400 \times 400$	Type 4	3.000e-02	2.094e+01
$600 \times 600$	Type 4	6.000e-02	5.456e + 01

Table 1: Execution time for different sizes and types of matrices

We can see that the power\_v11 algorithm is generally slower than the eigen function especially for the type 2 and 4 matrices.

## 1.2.

```
nb_it = 1;
norme = norm(beta*v - z, 2)/norm(beta,2);

while(norme > eps && nb_it < maxit)
    beta_old = beta;
    v = z/norm(z, 2);
    z = A*v;
    beta = (v'*z)/(v'*v);
    norme = abs(beta-beta_old)/abs(beta_old);
    nb_it = nb_it + 1;
end</pre>
```

Listing 1: Inner loop of the new algorithm

Matrix dimension	Matrix type	Exec. time for $power_v11$ , (s)	Exec. time for $power_v12$ , (s)
$200 \times 200$	Type 1	1.960e+00	3.200e-01
$400 \times 400$	Type 1	1.888e+01	2.660e+00
$600 \times 600$	Type 1	5.031e+01	7.070e+00
$200 \times 200$	Type 2	1.000e-02	1.000e-02
$400 \times 400$	Type 2	7.000e-02	1.000e-02
$600 \times 600$	Type 2	1.800e-01	4.000e-02
$200 \times 200$	Type 3	3.000e-02	1.000e-02
$400 \times 400$	Type 3	6.100e-01	1.100e-01
$600 \times 600$	Type 3	1.270e+00	2.600e-01

$200 \times 200$	Type 4	1.530e+00	2.900e-01
$400 \times 400$	Type 4	2.113e+01	3.060e+00
$600 \times 600$	Type 4	5.914 + e01	6.480e+00

We can see that the power\_v12 algorithm is globally faster than the power\_v11.

#### 1.3.

The main drawback of the deflated power method is the numerous matrix-vector products required to compute the eigenvectors as well as the fact that each iteration compute only one eigenvalue which can be slow if a lot of eigenvalues are desired.

#### 1.4.

If we apply Algorithm 1 to m vectors, there is no reason for the columns of V to converge to a base. Each vector will converge toward a different projection of the dominant eigenvalue.

#### 1.5.

In Algorithm 2, the matrix H is a smaller matrix, with dimension  $n \times m$ , therefore, even for larger matrices A, computing the spectral decomposition of H will not be computionally expensive.

## 2.

#### 2.1.

 $\Sigma_k$  is of size (k,k)

 $U_k$  is of size (q, k)

 $V_k$  is of size (p, k)

# 2.2.

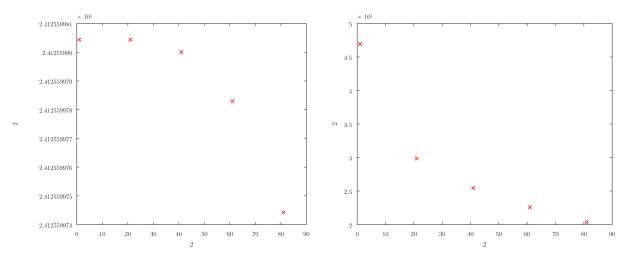
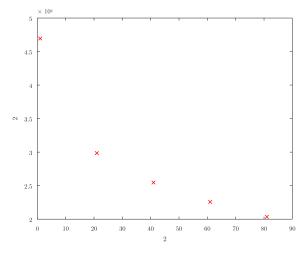


Figure 1: eig method

Figure 2: power method



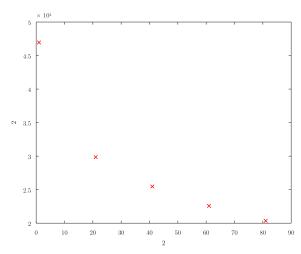


Figure 3: subspace\_iter0 method

 $Figure~4:~subspace\_iter1~method$ 

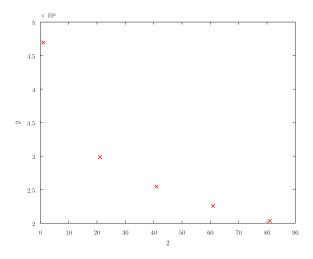


Figure 5:  $subspace\_iter2$  method