

November 2, 2023

Suppose we want to define the smallest Eigenvalue:

*Input:*  $A, v_0, \text{tol}, \text{matIter}$

```
while(error < tol && iter < matIter)
{ v1 = 1 / magnitude(y) * y
w = A * v1
 $\lambda_1 = v_1^t * w$ 
```

```
error = abs( $\lambda_0 - \lambda_1$ )
iter++;
 $\lambda_0 = \lambda_1$ 
```

```
y = v1
```

```
}
return  $\lambda_1$ , error;
```

**Finding the smallest Eigenvalue:** If  $\lambda_1$  is the largest eigenvalue of  $A$ , then  $\frac{1}{\lambda_1}$  is the smallest eigenvalue of  $A^{-1}$ . If  $\lambda_n$  is the smallest eigenvalue of  $A$ , then  $\frac{1}{\lambda_n}$  is the largest eigenvalue of  $A^{-1}$ .

We have the means to compute approximations of  $\lambda_1$  and  $\lambda_n$

- $\lambda_1$ : Use the power method
- $\lambda_n$ : Use the inverse power method

**properties of eigenvalues:**

if  $(\lambda, v)$  an eigen pair,  $v \neq 0$ . this means  $Av = \lambda v$ . Then for  $\mu \in \mathbb{R}$ ,  $(Av - \mu Iv) = (A - \mu I)v = \lambda v - \mu Iv = (\lambda - \mu)v$

$\rightarrow (\lambda - \mu)$  is an eigenvalue of  $(A - \mu I)$ .

So,  $(A - \mu I)v = (\lambda - \mu)v$

The idea is choose  $\mu$  close our eigenvalue. We can use inverse iteration to derive an approximation of  $\lambda - \mu$  and then add  $\mu$  back to get  $\lambda$ .

**Iteratively finding eigenvalues:**

$$v_0 = a_1 v_1 + a_2 v_2 + \dots + a_n v_n$$

$$Av_0 = a_1(\lambda_1 v_1) + a_2(\lambda_2 v_2) + \dots + a_n(\lambda_n v_n)$$

where  $\lambda_1 / > \lambda_2 \geq \dots \geq \lambda_{n-1} / > 0$

### Properties

$$1. \frac{Av_k}{\|v_k\|} \rightarrow v_1$$

$$2. \frac{v_k^t A v_k}{v_k^t v_k}$$

3. If  $\lambda$  is an eigenvalue of  $A$ , the  $\frac{1}{\lambda}$  is an eigenvalue of  $A^{-1}$

$$4. A\vec{v} = \lambda\vec{v} \rightarrow A\vec{v} - \mu\vec{v} = (\lambda - \mu)\vec{v} \rightarrow (A - \mu I)\vec{v}.$$

If  $\lambda$  is an eigenvalue of  $A$ , then  $\lambda - \mu$  is an eigenvalue of  $A - \mu I$

Using this approach allows us to utilize parallelism. This means we can use a tool like *openMP* to easily parallelize code written in C, C++, fortran, etc

### Lamczos Algorithm

$$Ax \rightarrow b$$

$$\begin{bmatrix} \sum_{j=1}^n a_{1,j}x_j \\ \sum_{j=1}^n a_{2,j}x_j \\ \dots \end{bmatrix} = \begin{bmatrix} b_1 \\ b_2 \\ \dots \\ b_n \end{bmatrix} =$$

$$x_1 \begin{bmatrix} a_{1,1} \\ a_{1,2} \\ \dots \\ a_{1,n} \end{bmatrix} + x_2 \begin{bmatrix} a_{2,1} \\ a_{2,2} \\ \dots \\ a_{2,n} \end{bmatrix} + \dots + x_n \begin{bmatrix} a_{n,1} \\ a_{n,2} \\ \dots \\ a_{n,n} \end{bmatrix}$$

```
for(int i = 0; i < n; i++){
```

```
double sum = 0.0;
```

```
for(int j = 0; j < n; j++){
```

```
sum += a[i][j] * x[j];
```

```
}
```

```
b[i] = sum;
```

```
}
```