

Power iteration

In mathematics, **power iteration** (also known as the *power method*) is an eigenvalue algorithm given a diagonalizable matrix **A**, the algorithm will produce a number **λ**, which is the greatest (in absolute value) eigenvalue of **A**, and a nonzero vector **v**, the corresponding eigenvector of **λ**, such that **Av = λv**. The algorithm is also known as the von Mises iteration^[1]

Power iteration is a very simple algorithm, but it may converge slowly. It does not compute a matrix decomposition and hence it can be used when **A** is a very large sparse matrix.

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The method

The power iteration algorithm starts with a vector **b**₀, which may be an approximation to the dominant eigenvector or a random vector. The method is described by the recurrence relation

$$b_{k+1} = \frac{Ab_k}{\|Ab_k\|}$$

So, at every iteration, the vector **b**_k is multiplied by the matrix **A** and normalized.

If we assume **A** has an eigenvalue that is strictly greater in magnitude than its other eigenvalues and the starting vector **b**₀ has a nonzero component in the direction of an eigenvector associated with the dominant eigenvalue, then a subsequence (**b**_k) converges to an eigenvector associated with the dominant eigenvalue.

Without the two assumptions above, the sequence (**b**_k) does not necessarily converge. In this sequence,

$$b_k = e^{i\phi_k} v_1 + r_k,$$

where **v**₁ is an eigenvector associated with the dominant eigenvalue, and **||r**_k**||** → 0. The presence of the term **e**^{*iφ_k*} implies that (**b**_k) does not converge unless **e**^{*iφ_k*} = 1. Under the two assumptions listed above, the sequence (**μ**_k) defined by

$$\mu_k = \frac{b_k^* Ab_k}{b_k^* b_k}$$

converges to the dominant eigenvalue

One may compute this with the following algorithm (shown in Python with NumPy):

```
#!/usr/bin/python
import numpy as np
```

```
def power_iteration(A, num_simulations):
    # Ideally choose a random vector
    # To decrease the chance that our vector
    # Is orthogonal to the eigenvector
    b_k = np.random.rand(A.shape[0])

    for _ in range(num_simulations):
        # calculate the matrix-by-vector product Ab
        b_k1 = np.dot(A, b_k)

        # calculate the norm
        b_k1_norm = np.linalg.norm(b_k1)

        # re normalize the vector
        b_k = b_k1 / b_k1_norm

    return b_k

power_iteration(np.array([[0.5, 0.5], [0.2, 0.8]]), 10)
```

The vector \mathbf{b}_k to an associated eigenvector. Ideally, one should use the Rayleigh quotient in order to get the associated eigenvalue.

This algorithm is the one used to calculate such things as the Google PageRank.

The method can also be used to calculate the spectral radius (the largest eigenvalue of a matrix) by computing the Rayleigh quotient

$$\frac{\mathbf{b}_k^\top \mathbf{A} \mathbf{b}_k}{\mathbf{b}_k^\top \mathbf{b}_k} = \frac{\mathbf{b}_{k+1}^\top \mathbf{b}_k}{\mathbf{b}_k^\top \mathbf{b}_k}.$$

Analysis

Let \mathbf{A} be decomposed into its Jordan canonical form: $\mathbf{A} = \mathbf{V} \mathbf{J} \mathbf{V}^{-1}$, where the first column of \mathbf{V} is an eigenvector of \mathbf{A} corresponding to the dominant eigenvalue λ_1 . Since the dominant eigenvalue of \mathbf{A} is unique, the first Jordan block of \mathbf{J} is the 1×1 matrix $[\lambda_1]$, where λ_1 is the largest eigenvalue of \mathbf{A} in magnitude. The starting vector \mathbf{b}_0 can be written as a linear combination of the columns of \mathbf{V} : $\mathbf{b}_0 = c_1 \mathbf{v}_1 + c_2 \mathbf{v}_2 + \dots + c_n \mathbf{v}_n$. By assumption, \mathbf{b}_0 has a nonzero component in the direction of the dominant eigenvalue, so $c_1 \neq 0$.

The computationally useful recurrence relation for \mathbf{b}_{k+1} can be rewritten as: $\mathbf{b}_{k+1} = \frac{\mathbf{A} \mathbf{b}_k}{\|\mathbf{A} \mathbf{b}_k\|} = \frac{\mathbf{A}^{k+1} \mathbf{b}_0}{\|\mathbf{A}^{k+1} \mathbf{b}_0\|}$, where the expression:

$\frac{\mathbf{A}^{k+1} \mathbf{b}_0}{\|\mathbf{A}^{k+1} \mathbf{b}_0\|}$ is more amenable to the following analysis.

$$\begin{aligned} \mathbf{b}_k &= \frac{\mathbf{A}^k \mathbf{b}_0}{\|\mathbf{A}^k \mathbf{b}_0\|} \\ &= \frac{(\mathbf{V} \mathbf{J} \mathbf{V}^{-1})^k \mathbf{b}_0}{\|(\mathbf{V} \mathbf{J} \mathbf{V}^{-1})^k \mathbf{b}_0\|} \\ &= \frac{\mathbf{V} \mathbf{J}^k \mathbf{V}^{-1} \mathbf{b}_0}{\|\mathbf{V} \mathbf{J}^k \mathbf{V}^{-1} \mathbf{b}_0\|} \\ &= \frac{\mathbf{V} \mathbf{J}^k \mathbf{V}^{-1} (c_1 \mathbf{v}_1 + c_2 \mathbf{v}_2 + \dots + c_n \mathbf{v}_n)}{\|\mathbf{V} \mathbf{J}^k \mathbf{V}^{-1} (c_1 \mathbf{v}_1 + c_2 \mathbf{v}_2 + \dots + c_n \mathbf{v}_n)\|} \\ &= \frac{\mathbf{V} \mathbf{J}^k (c_1 \mathbf{e}_1 + c_2 \mathbf{e}_2 + \dots + c_n \mathbf{e}_n)}{\|\mathbf{V} \mathbf{J}^k (c_1 \mathbf{e}_1 + c_2 \mathbf{e}_2 + \dots + c_n \mathbf{e}_n)\|} \\ &= \left(\frac{\lambda_1}{|\lambda_1|} \right)^k \frac{c_1}{|c_1|} \frac{v_1 + \frac{1}{c_1} V \left(\frac{1}{\lambda_1} J \right)^k (c_2 \mathbf{e}_2 + \dots + c_n \mathbf{e}_n)}{\|v_1 + \frac{1}{c_1} V \left(\frac{1}{\lambda_1} J \right)^k (c_2 \mathbf{e}_2 + \dots + c_n \mathbf{e}_n)\|} \end{aligned}$$

The expression above simplifies as $k \rightarrow \infty$

$$\left(\frac{1}{\lambda_1} J\right)^k = \begin{bmatrix} [1] & & & \\ & \left(\frac{1}{\lambda_1} J_2\right)^k & & \\ & & \ddots & \\ & & & \left(\frac{1}{\lambda_1} J_m\right)^k \end{bmatrix} \rightarrow \begin{bmatrix} 1 & & & \\ & 0 & & \\ & & \ddots & \\ & & & 0 \end{bmatrix} \text{ as } k \rightarrow \infty.$$

The limit follows from the fact that the eigenvalue of $\frac{1}{\lambda_1} J_i$ is less than 1 in magnitude, so $\left(\frac{1}{\lambda_1} J_i\right)^k \rightarrow 0$ as $k \rightarrow \infty$

It follows that:

$$\frac{1}{c_1} V \left(\frac{1}{\lambda_1} J\right)^k (c_2 e_2 + \dots + c_n e_n) \rightarrow 0 \text{ as } k \rightarrow \infty$$

Using this fact, b_k can be written in a form that emphasizes its relationship with v_1 when k is large:

$$b_k = \left(\frac{\lambda_1}{|\lambda_1|}\right)^k \frac{c_1}{|c_1|} \frac{v_1 + \frac{1}{c_1} V \left(\frac{1}{\lambda_1} J\right)^k (c_2 e_2 + \dots + c_n e_n)}{\|v_1 + \frac{1}{c_1} V \left(\frac{1}{\lambda_1} J\right)^k (c_2 e_2 + \dots + c_n e_n)\|} = e^{i\phi_k} \frac{c_1}{|c_1|} \frac{v_1}{\|v_1\|} + r_k \text{ where } e^{i\phi_k} = (\lambda_1/|\lambda_1|)^k \text{ and } \|r_k\| \rightarrow 0 \text{ as } k \rightarrow \infty$$

The sequence (b_k) is bounded, so it contains a convergent subsequence. Note that the eigenvector corresponding to the dominant eigenvalue is only unique up to a scalar, so although the sequence (b_k) may not converge, b_k is nearly an eigenvector of A for large k .

Alternatively, if A is diagonalizable, then the following proof yields the same result

Let $\lambda_1, \lambda_2, \dots, \lambda_m$ be the m eigenvalues (counted with multiplicity) of A and let v_1, v_2, \dots, v_m be the corresponding eigenvectors. Suppose that λ_1 is the dominant eigenvalue, so that $|\lambda_1| > |\lambda_j|$ for $j > 1$.

The initial vector b_0 can be written:

$$b_0 = c_1 v_1 + c_2 v_2 + \dots + c_m v_m.$$

If b_0 is chosen randomly (with uniform probability), then $c_1 \neq 0$ with probability 1. Now,

$$\begin{aligned} A^k b_0 &= c_1 A^k v_1 + c_2 A^k v_2 + \dots + c_m A^k v_m \\ &= c_1 \lambda_1^k v_1 + c_2 \lambda_2^k v_2 + \dots + c_m \lambda_m^k v_m \\ &= c_1 \lambda_1^k \left(v_1 + \frac{c_2}{c_1} \left(\frac{\lambda_2}{\lambda_1}\right)^k v_2 + \dots + \frac{c_m}{c_1} \left(\frac{\lambda_m}{\lambda_1}\right)^k v_m \right). \end{aligned}$$

The expression within parentheses converges to v_1 because $|\lambda_j/\lambda_1| < 1$ for $j > 1$. On the other hand, we have

$$b_k = \frac{A^k b_0}{\|A^k b_0\|}.$$

Therefore, b_k converges to (a multiple of) the eigenvector v_1 . The convergence is geometric, with ratio

$$\left| \frac{\lambda_2}{\lambda_1} \right|,$$

where λ_2 denotes the second dominant eigenvalue. Thus, the method converges slowly if there is an eigenvalue close in magnitude to the dominant eigenvalue.

Applications

Although the power iteration method approximates only one eigenvalue of a matrix, it remains useful for certain computational problems. For instance, Google uses it to calculate the PageRank of documents in their search engine,^[2] and Twitter uses it to show users recommendations of who to follow.^[3] For matrices that are well-conditioned and as sparse as the web matrix, the power

iteration method can be more efficient than other methods of finding the dominant eigenvector

Some of the more advanced eigenvalue algorithms can be understood as variations of the power iteration. For instance, the inverse iteration method applies power iteration to the matrix \mathbf{A}^{-1} . Other algorithms look at the whole subspace generated by the vectors \mathbf{b}_k . This subspace is known as the Krylov subspace. It can be computed by Arnoldi iteration or Lanczos iteration.

See also

- Rayleigh quotient iteration
- Inverse iteration

References

1. Richard von Mises and H. Pollaczek-Geiringer, *Praktische Verfahren der Gleichungsauflösung*, ZAMM - Zeitschrift für Angewandte Mathematik und Mechanik 9, 152-164 (1929).
2. Ipsen, Ilse, and Rebecca M. Wills (5–8 May 2005). "7th IMACS International Symposium on Iterative Methods in Scientific Computing" (http://www4.ncsu.edu/~ipsen/ps/slides_imacs.pdf) (PDF). Fields Institute, Toronto, Canada.
3. Pankaj Gupta, Ashish Goel, Jimmy Lin, Aneesh Sharma, Dong Wang, and Reza Bosagh Zadeh **WTF: The who-to-follow system at Twitter** (<http://dl.acm.org/citation.cfm?id=2488433>) Proceedings of the 22nd international conference on World Wide Web

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This page was last edited on 8 January 2018, at 21:33.

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