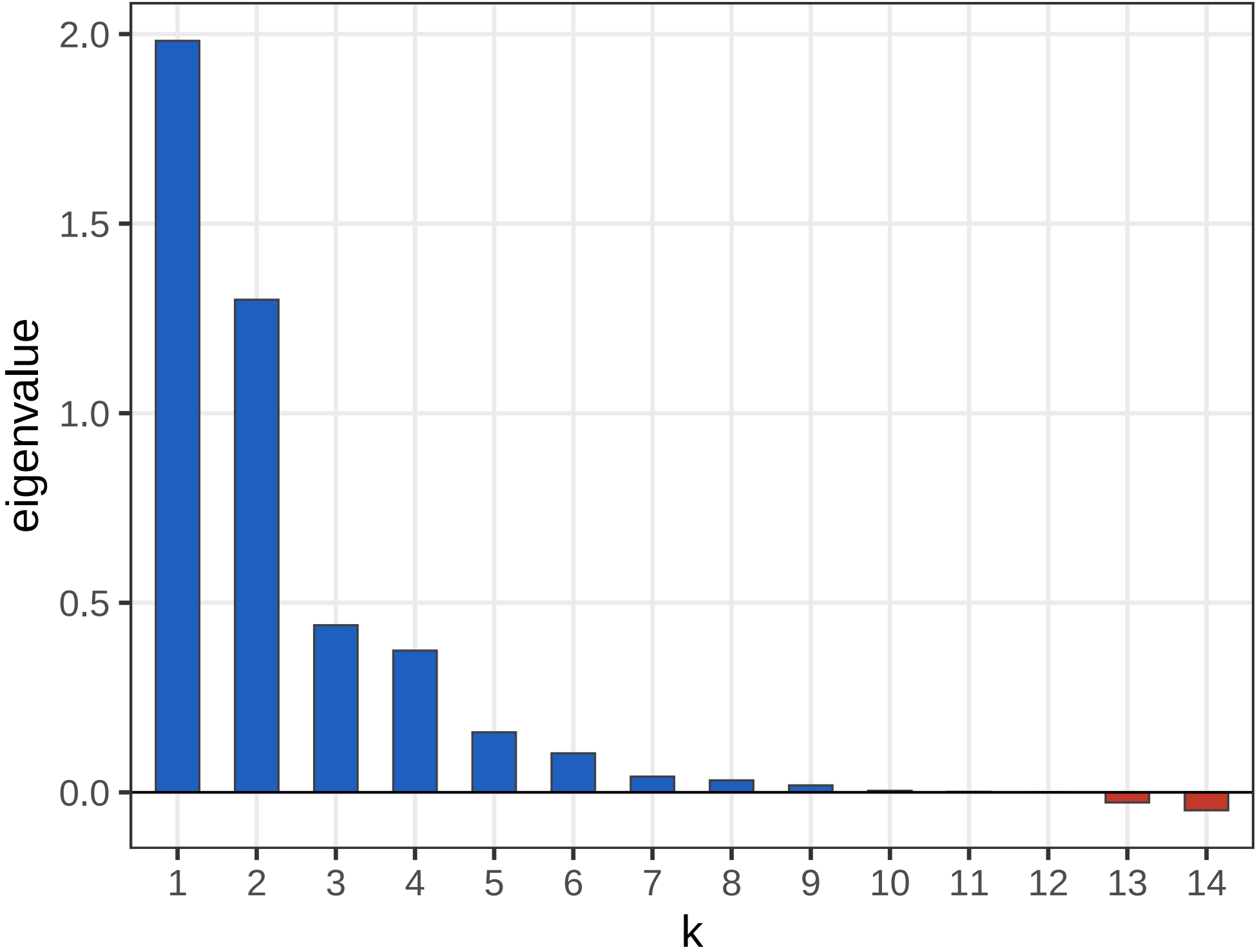

2.0
1.5
1.0
0.5
0.0
eigenvalue
1
2
3
4
5
6
7
8
9
10
11
12
13
14
k