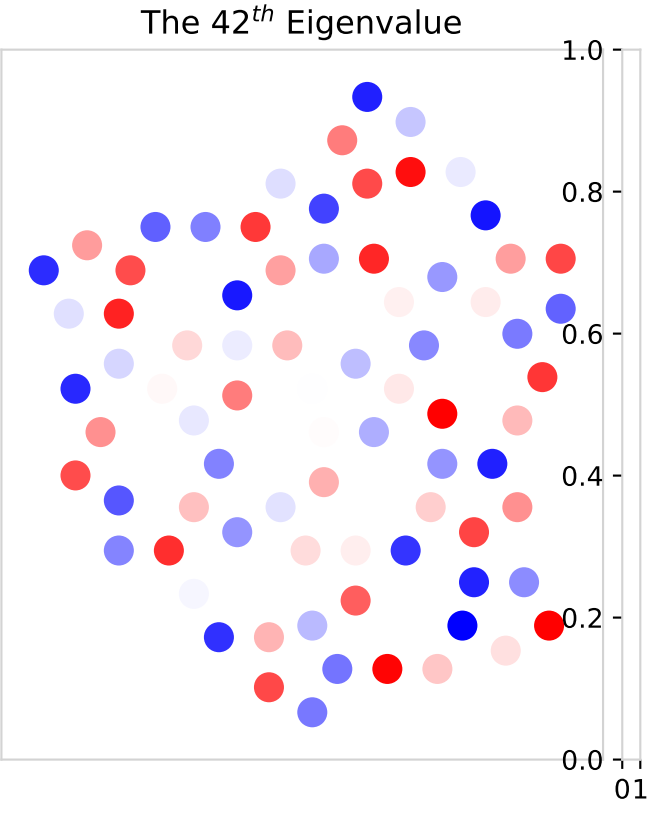
The 42th Eigenvalue
1.0
0.8
0.6
0.4
0.2
0.0
01