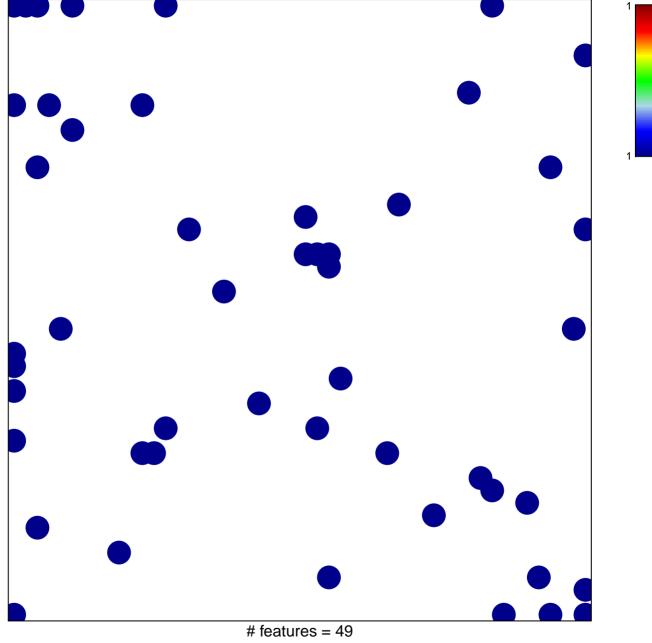
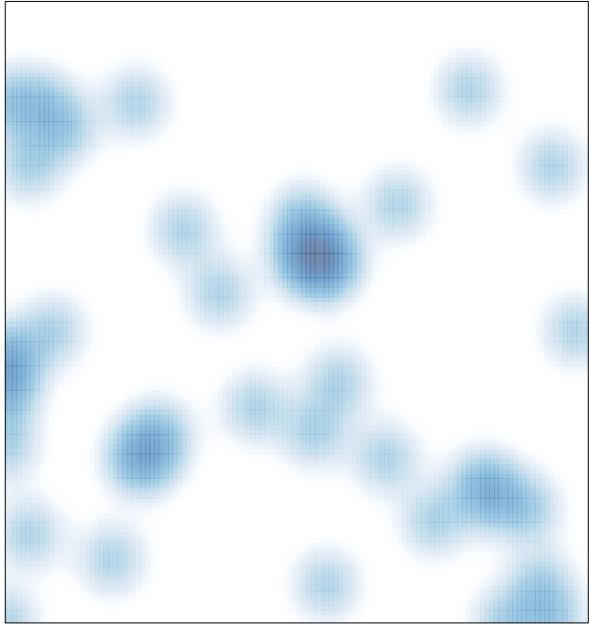
drug binding



chi-square p = 0.8

drug binding



features = 49, max = 1