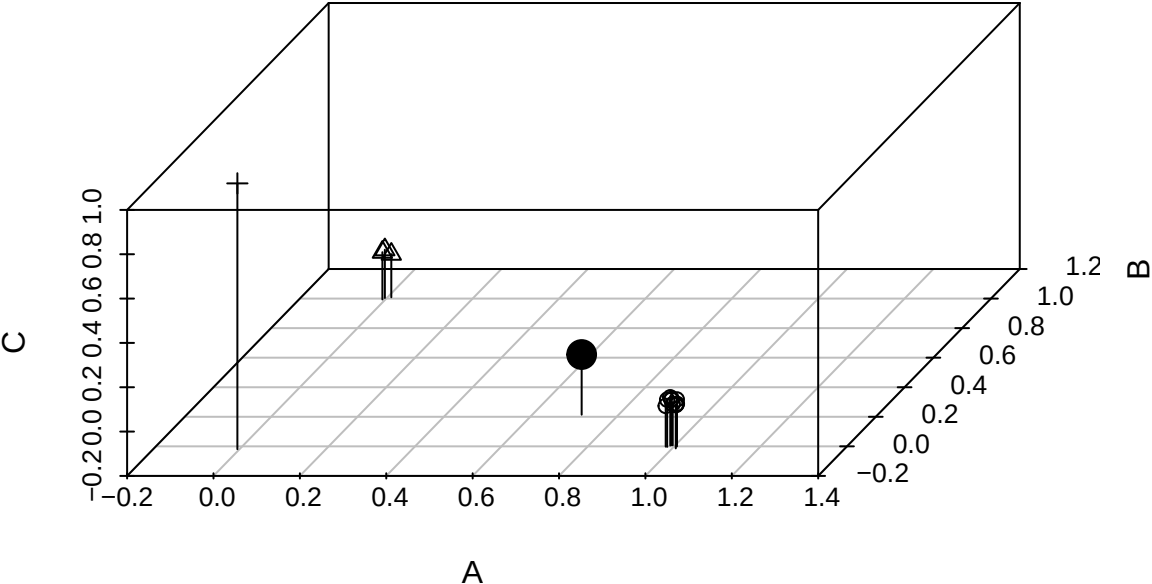
C
−0.2
0.0
0.2
0.4
0.6
0.8
1.0
A
−0.2
0.0
0.2
0.4
0.6
0.8
1.0
1.2
1.4
B
−0.2
0.0
0.2
0.4
0.6
0.8
1.0
1.2