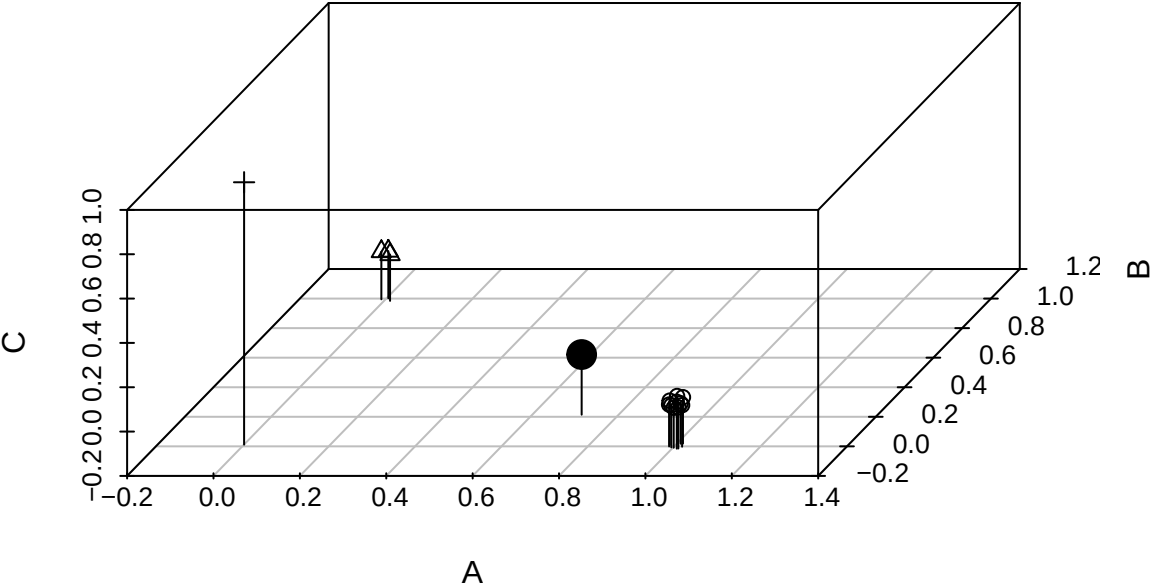
C
−0.2 0.0 0.2 0.4 0.6 0.8 1.0
A
−0.2 0.0 0.2 0.4 0.6 0.8 1.0 1.2 1.4
B
−0.2 0.0 0.2 0.4 0.6 0.8 1.0 1.2