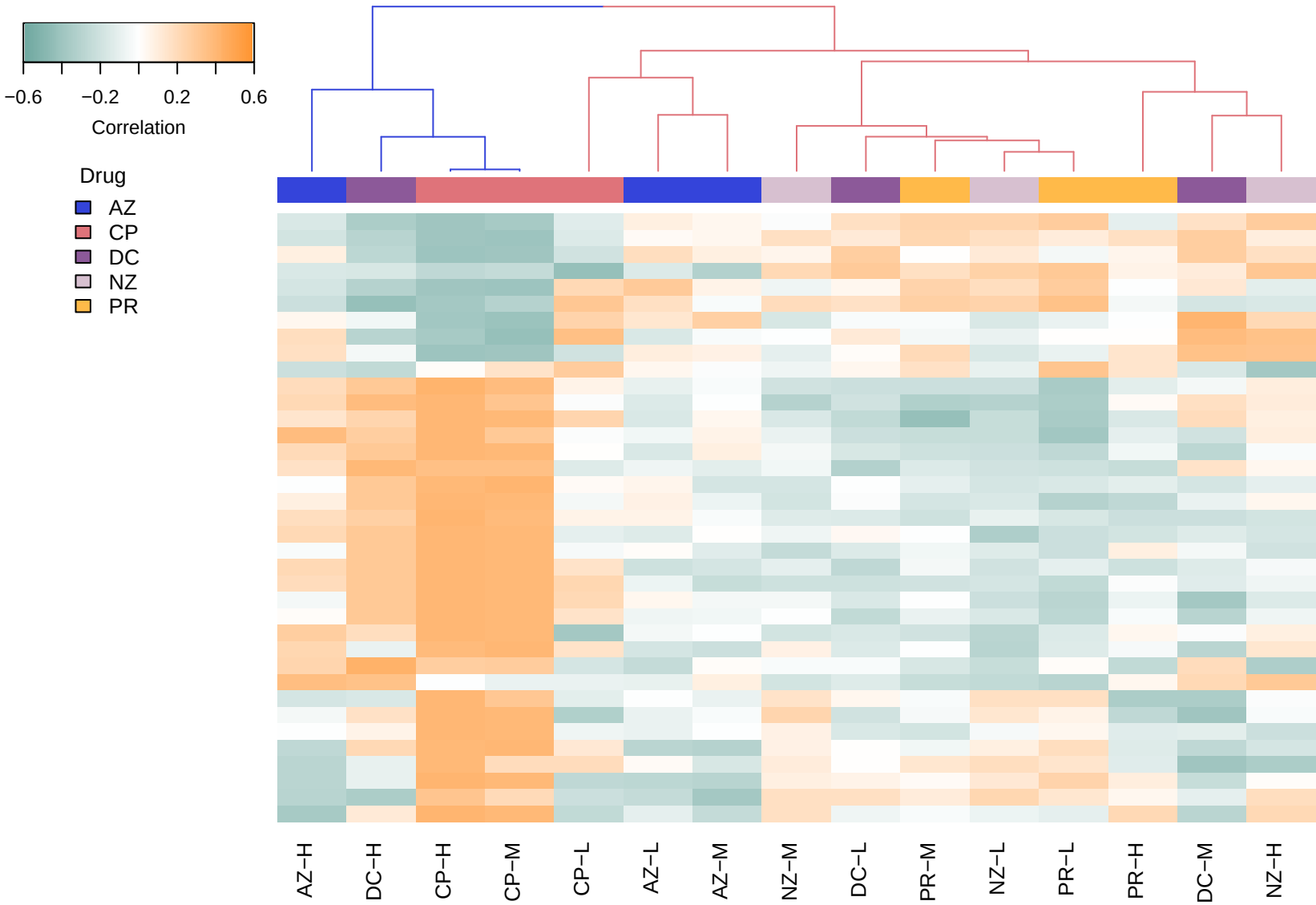
−0.6
−0.2
0.2
0.6
Correlation
Drug
AZ
CP
DC
NZ
PR
AZ−H
DC−H
CP−H
CP−M
CP−L
AZ−L
AZ−M
NZ−M
DC−L
PR−M
NZ−L
PR−L
PR−H
DC−M
NZ−H