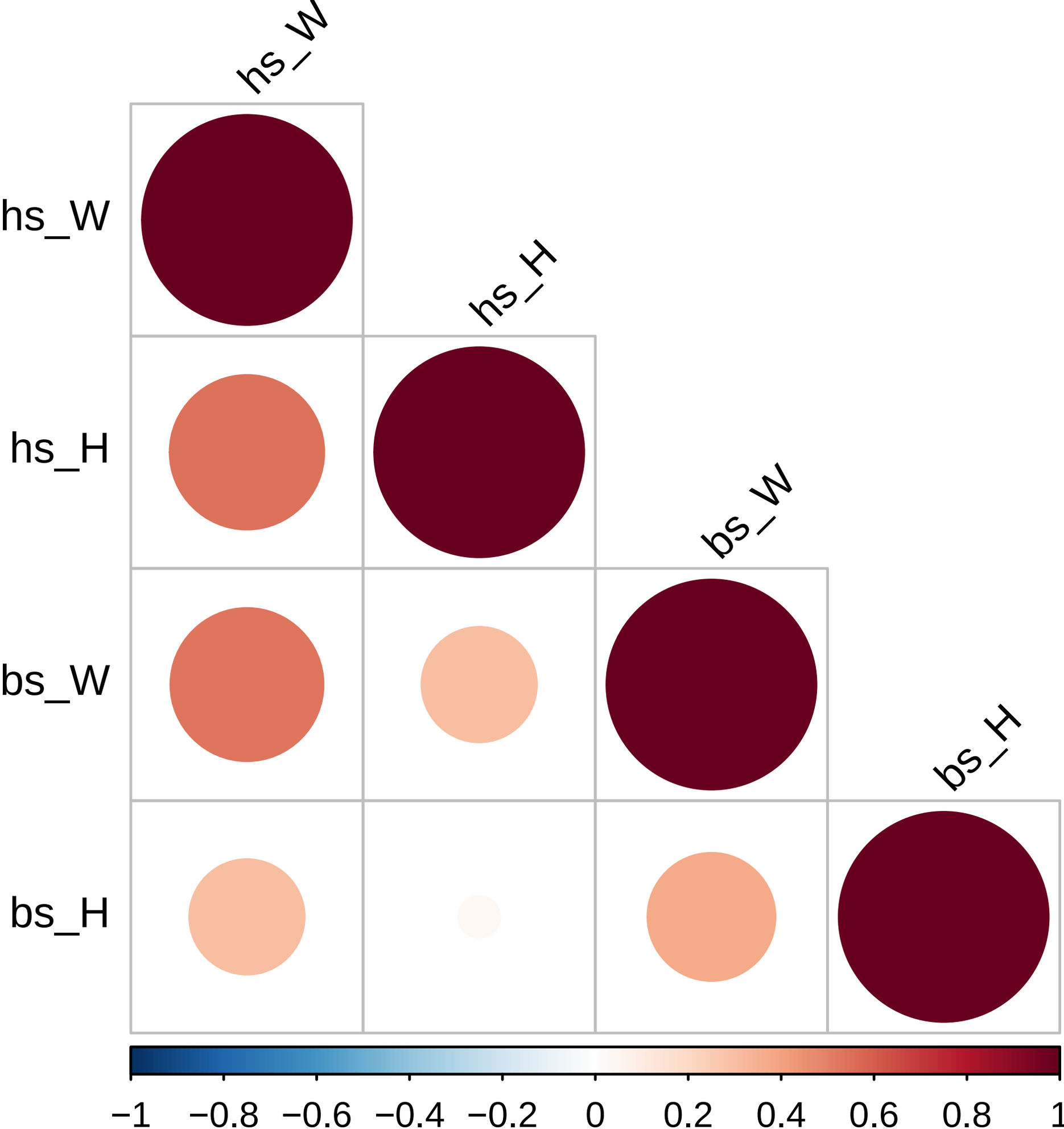
hs_W
hs_H
bs_W
bs_H
−1
−0.8
−0.6
−0.4
−0.2
0
0.2
0.4
0.6
0.8
1