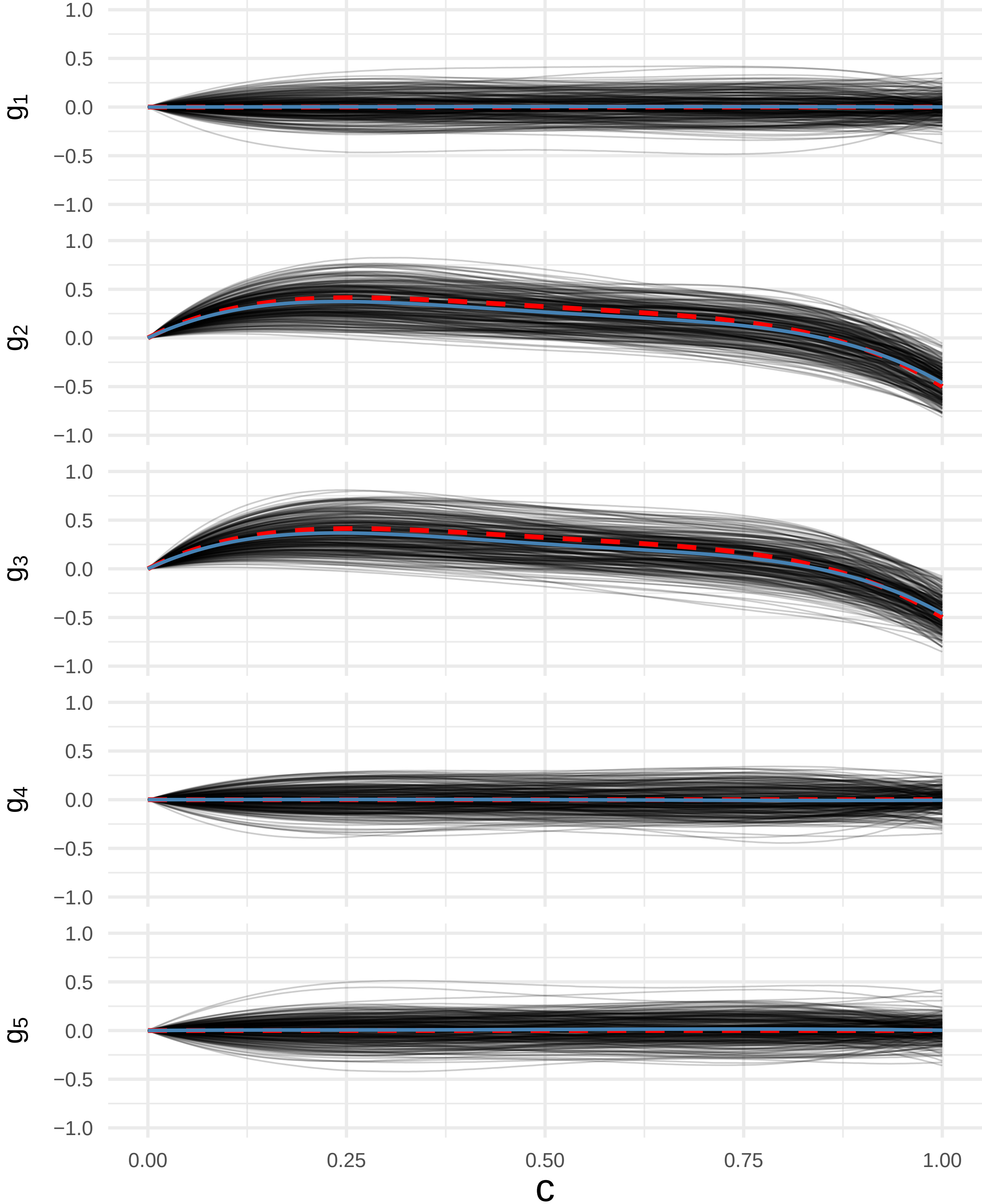
g_1
g_2
g_3
g_4
g_5
1.0
0.5
0.0
−0.5
−1.0
0.00
0.25
0.50
0.75
1.00
c