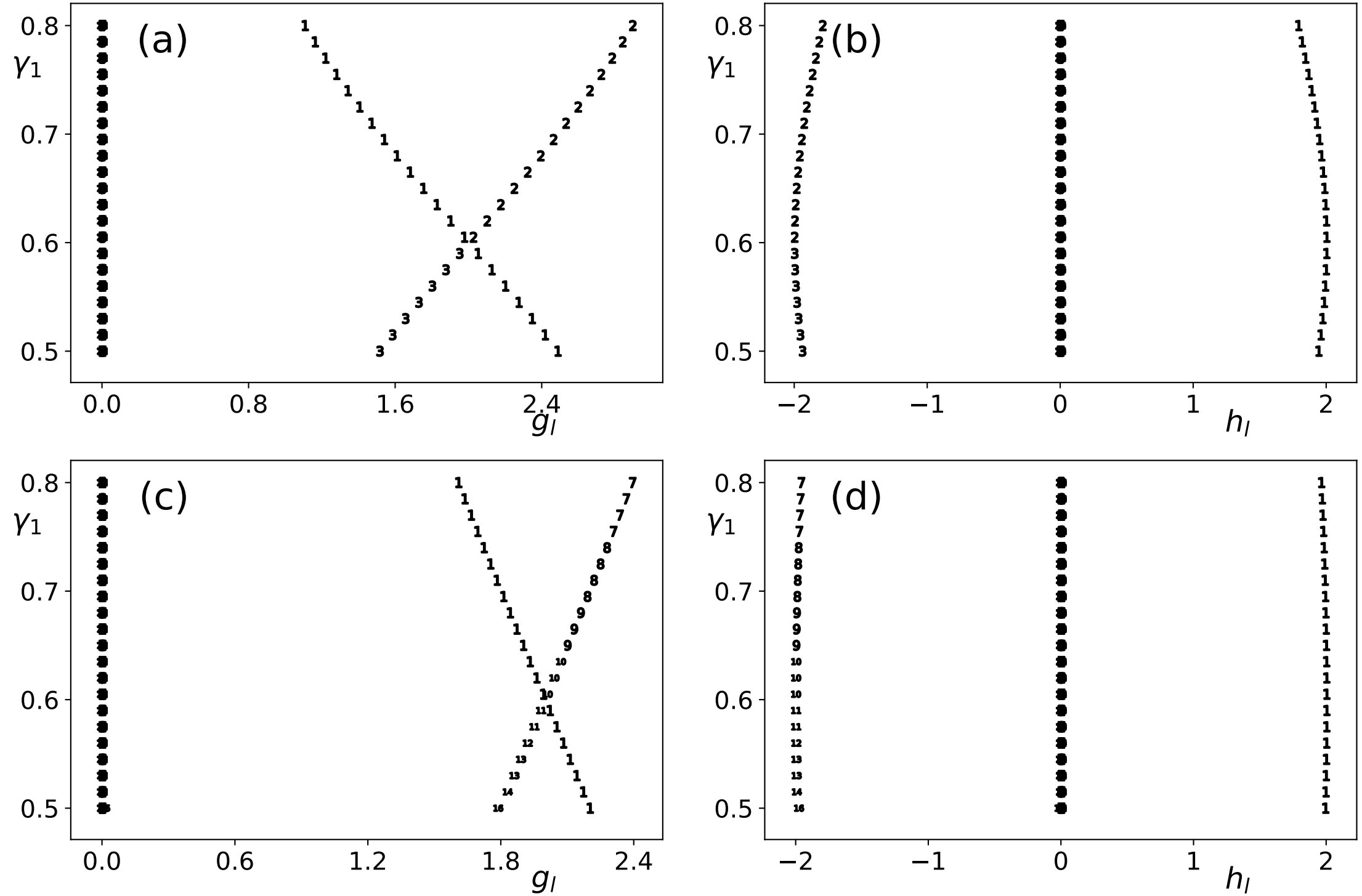
(a)
(b)
(c)
(d)
γ_1
g_l
h_l
0.8
0.7
0.6
0.5
0.0
0.8
1.6
2.4
0.6
1.2
1.8
−2
−1
0
1
2