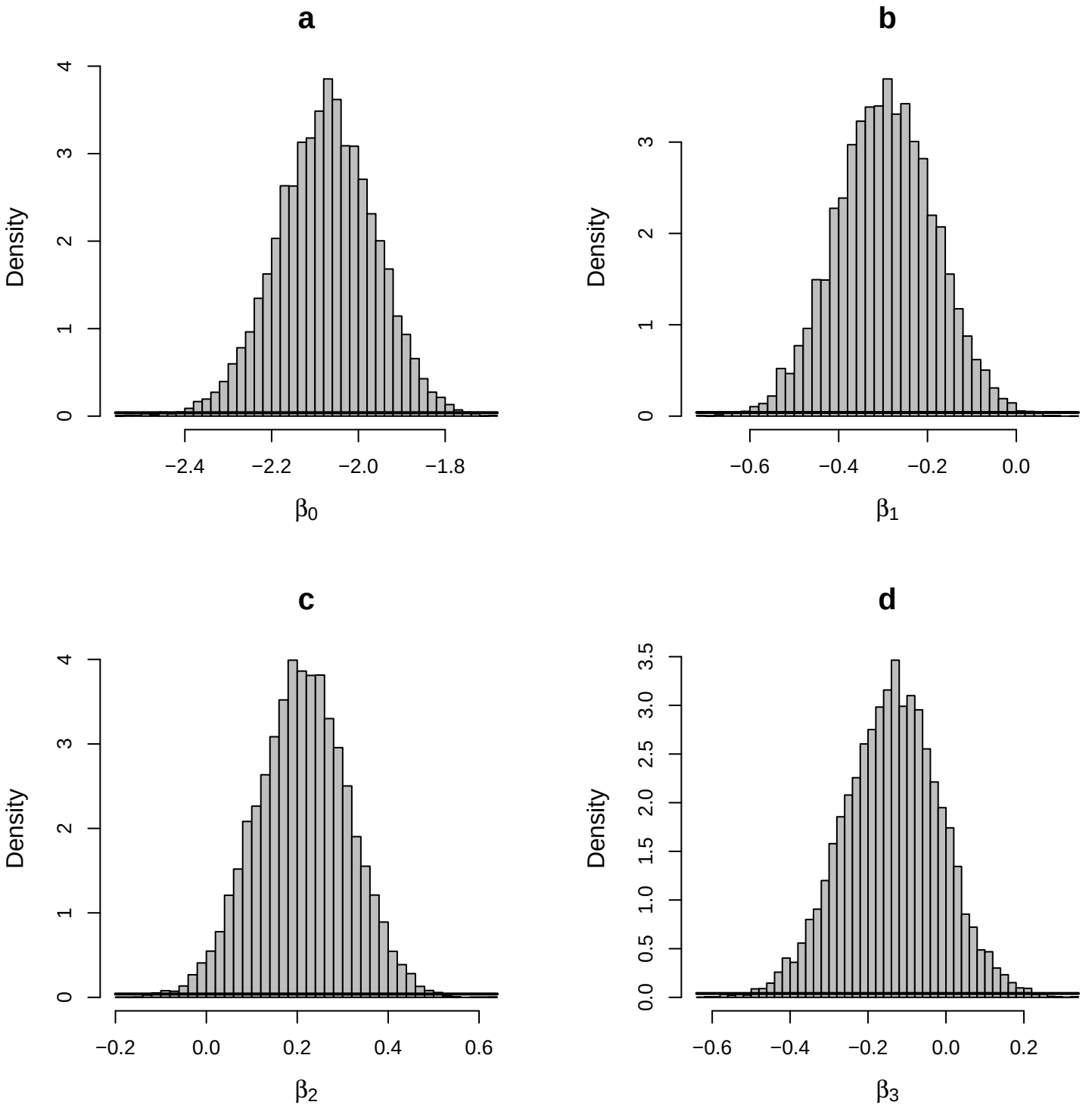
a
Density
0
1
2
3
4
−2.4
−2.2
−2.0
−1.8
β0
b
Density
0
1
2
3
−0.6
−0.4
−0.2
0.0
β1
c
Density
0
1
2
3
4
−0.2
0.0
0.2
0.4
0.6
β2
d
Density
0.0
0.5
1.0
1.5
2.0
2.5
3.0
3.5
−0.6
−0.4
−0.2
0.0
0.2
β3