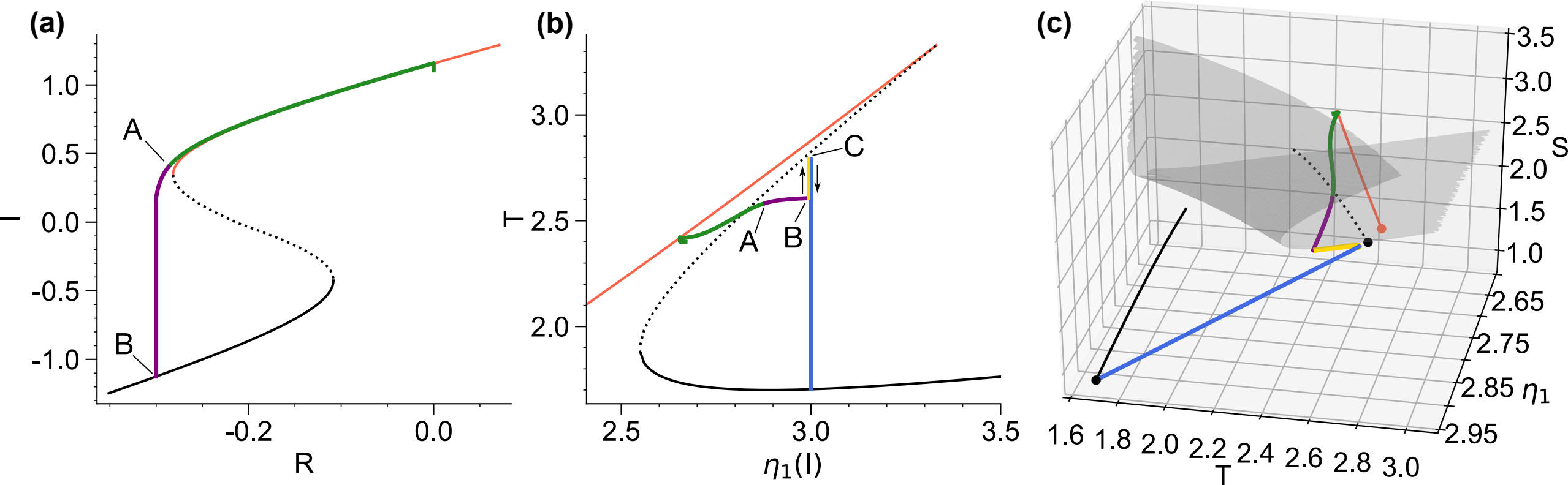
(a)
1.0
0.5
0.0
-0.5
-1.0
A
B
-0.2
0.0
R
(b)
3.0
2.5
2.0
C
A
B
2.5
3.0
3.5
η_1(I)
(c)
3.5
3.0
2.5
2.0
1.5
1.0
S
2.65
2.75
2.85
η_1
2.95
1.6 1.8 2.0 2.2 2.4 2.6 2.8 3.0
T