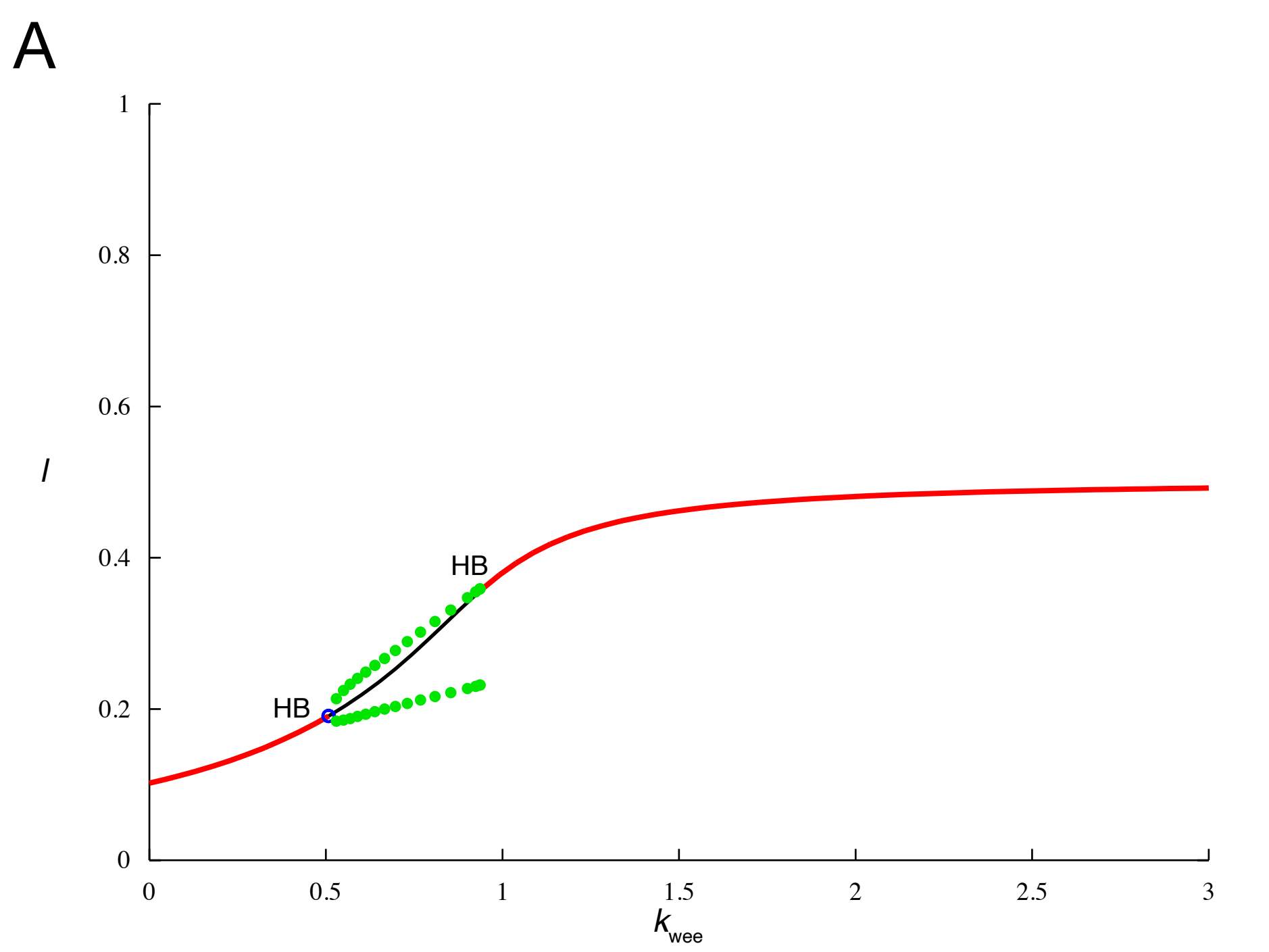
A
1
0.8
0.6
0.4
0.2
0
I
HB
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0
0.5
1
1.5
2
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3
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