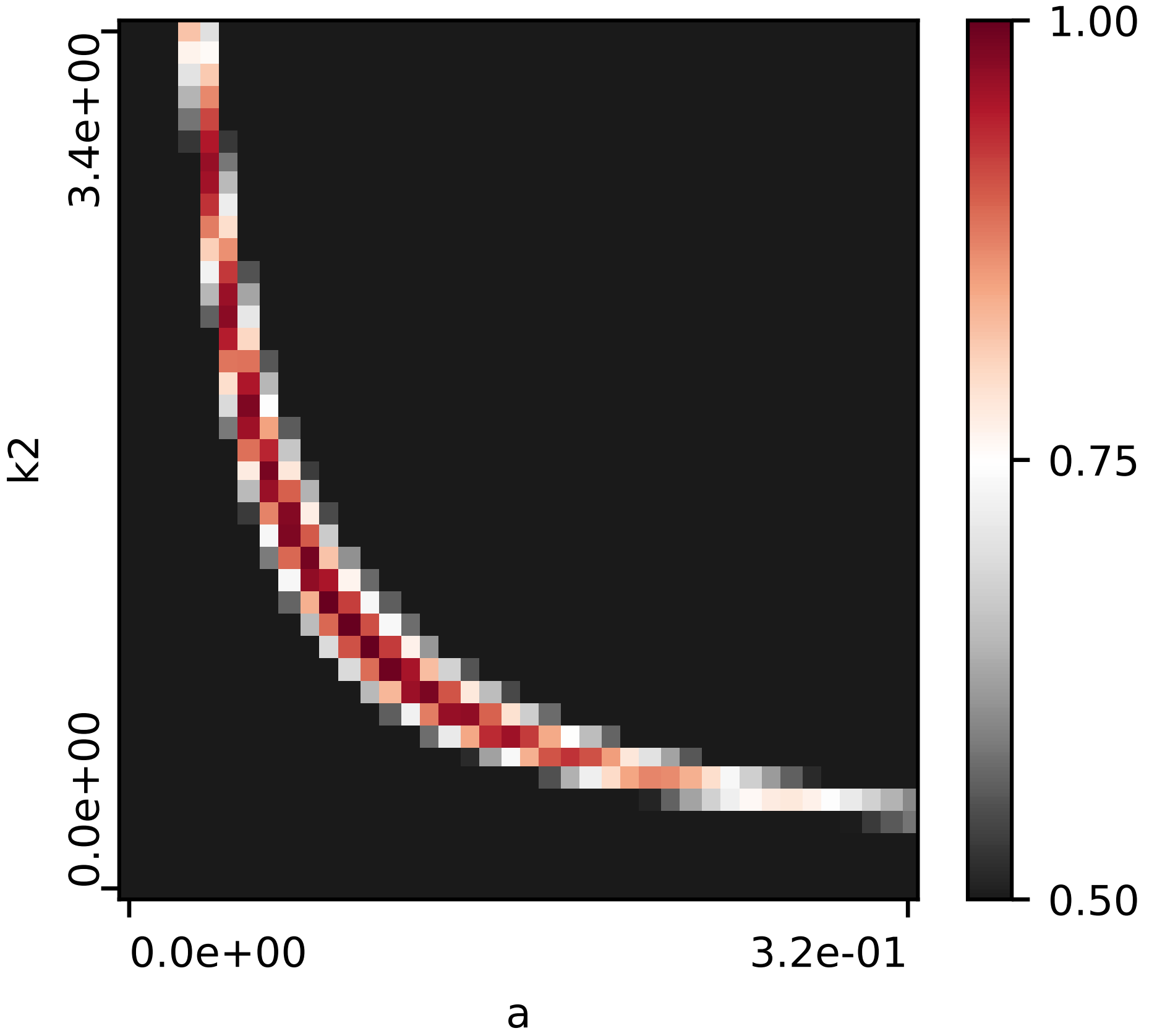

1.00
0.75
0.50
3.4e+00
0.0e+00
k2
0.0e+00
3.2e-01
a