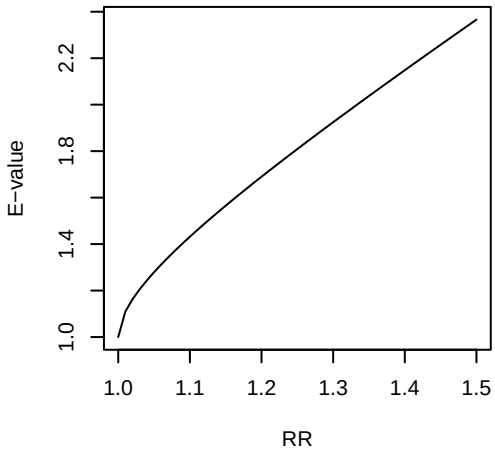
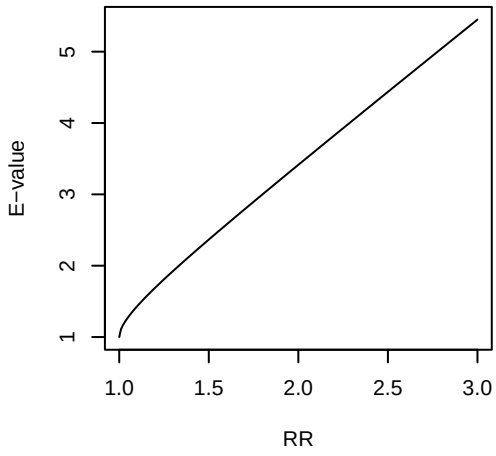
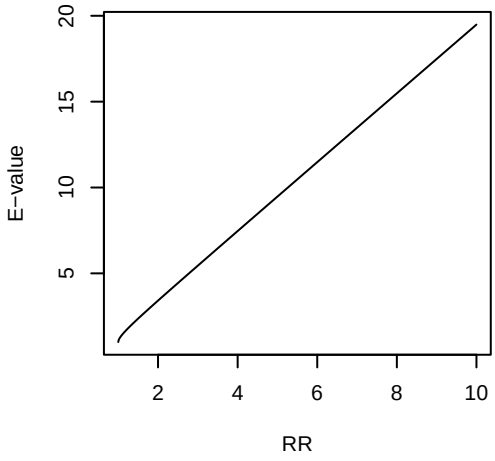

E−value
1.0
1.4
1.8
2.2
1.0
1.1
1.2
1.3
1.4
1.5
RR


E−value
1
2
3
4
5
1.0
1.5
2.0
2.5
3.0
RR


E−value
5
10
15
20
2
4
6
8
10
RR