(a)
Atomistic (pvpH)
< No. of H-bonds >
0
2
4
6
8
0
100
200
300
400
500
Simulation time [ns]

(b)
Atomistic (pvpH)
< R_ee > [d]
0
2
4
6
8
10
0
100
200
300
400
500
Simulation time [ns]