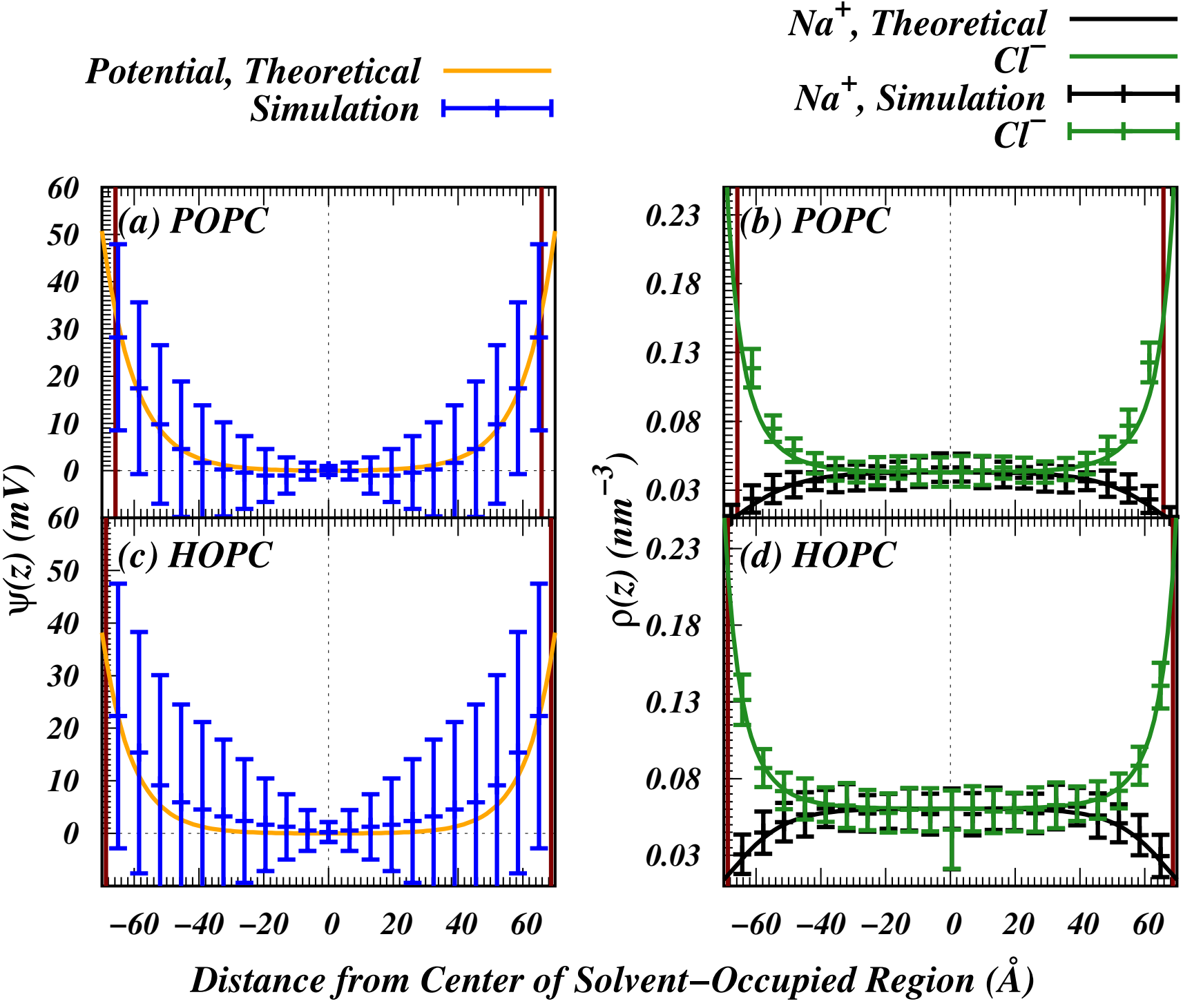
Potential, Theoretical
Simulation
Na+, Theoretical
Cl−
Na+, Simulation
Cl−
(a) POPC
(b) POPC
(c) HOPC
(d) HOPC
ψ(z) (mV)
ρ(z) (nm−3)
60
50
40
30
20
10
0
0.23
0.18
0.13
0.08
0.03
−60 −40 −20 0 20 40 60
Distance from Center of Solvent−Occupied Region (Å)