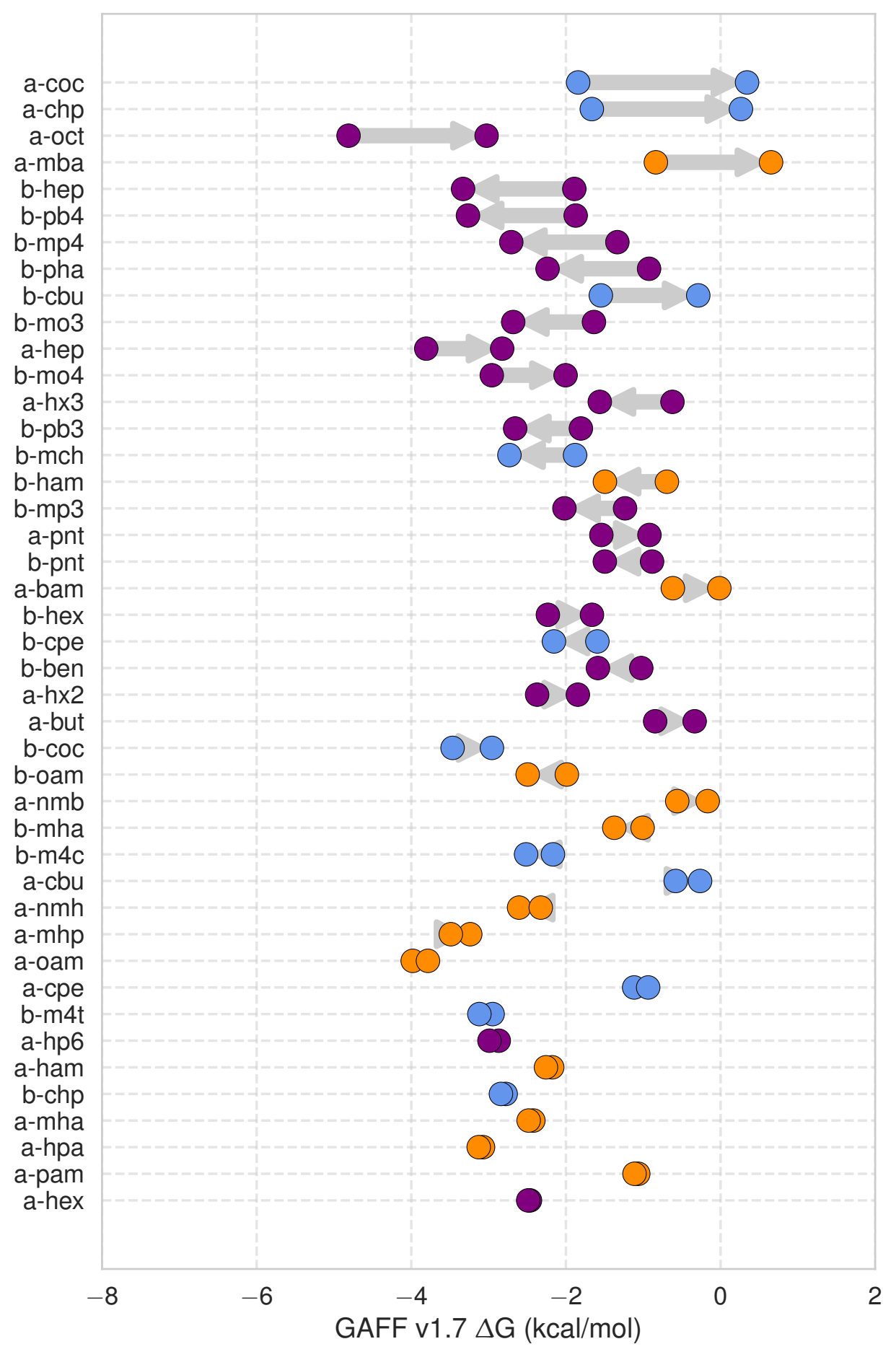
a-coc
a-chp
a-oct
a-mba
b-hep
b-pb4
b-mp4
b-pha
b-cbu
b-mo3
a-hep
b-mo4
a-hx3
b-pb3
b-mch
b-ham
b-mp3
a-pnt
b-pnt
a-bam
b-hex
b-cpe
b-ben
a-hx2
a-but
b-coc
b-oam
a-nmb
b-mha
b-m4c
a-cbu
a-nmh
a-mhp
a-oam
a-cpe
b-m4t
a-hp6
a-ham
b-chp
a-mha
a-hpa
a-pam
a-hex
−8
−6
−4
−2
0
2
GAFF v1.7 ΔG (kcal/mol)