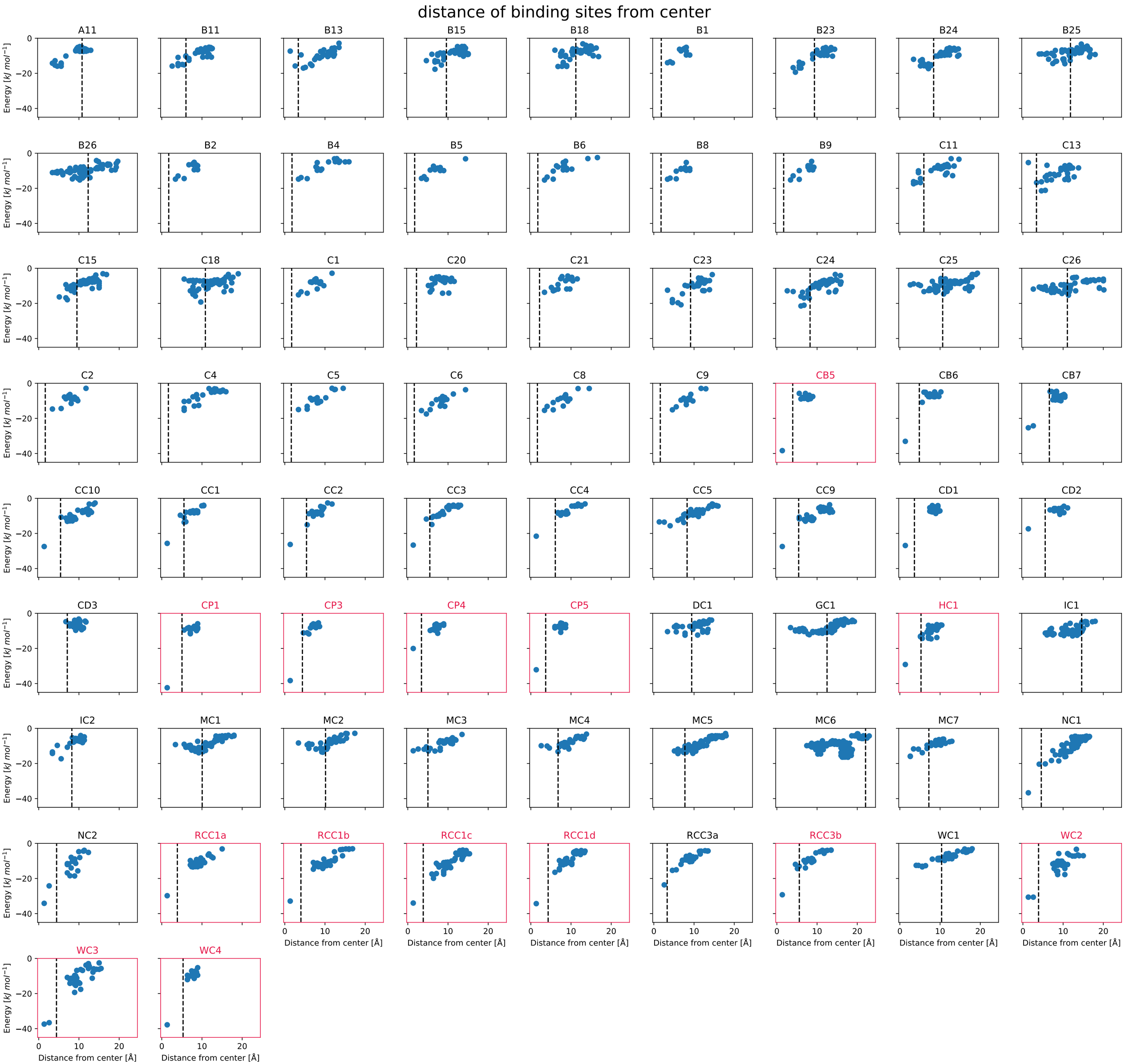
distance of binding sites from center
A11
B11
B13
B15
B18
B1
B23
B24
B25
B26
B2
B4
B5
B6
B8
B9
C11
C13
C15
C18
C1
C20
C21
C23
C24
C25
C26
C2
C4
C5
C6
C8
C9
CB5
CB6
CB7
CC10
CC1
CC2
CC3
CC4
CC5
CC9
CD1
CD2
CD3
CP1
CP3
CP4
CP5
DC1
GC1
HC1
IC1
IC2
MC1
MC2
MC3
MC4
MC5
MC6
MC7
NC1
NC2
RCC1a
RCC1b
RCC1c
RCC1d
RCC3a
RCC3b
WC1
WC2
WC3
WC4
Energy [kJ mol⁻¹]
0
−20
−40
0
10
20
Distance from center [Å]