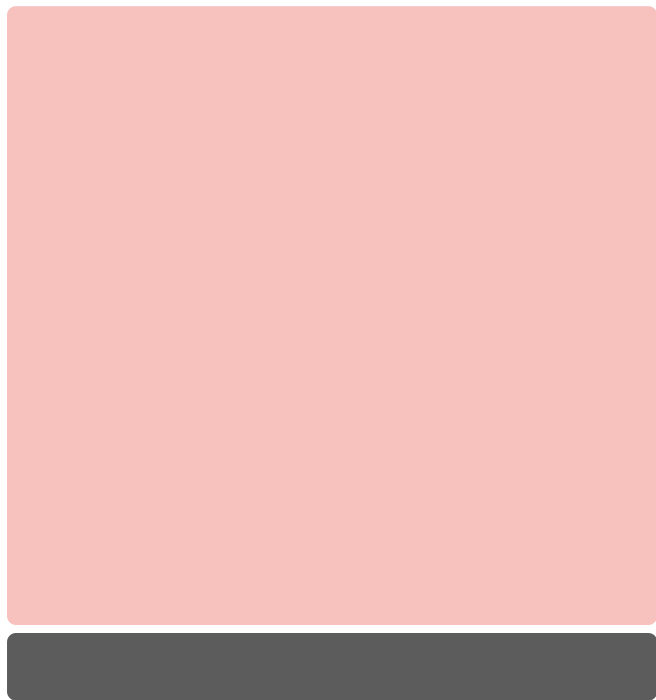


# (k-MDCM + CC) Energy Function + Relative Errors

Coulomb term



Interaction Energy



Pair potential term

