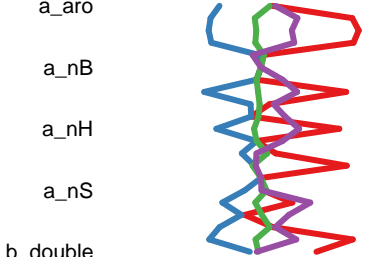


(EII,4,6014) Adjacency, distance matrix, connectivity, Kappa shape in



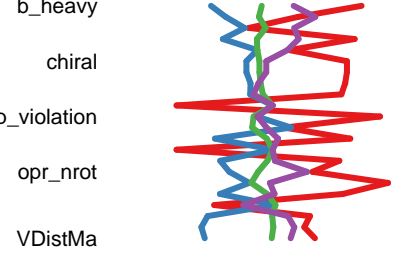
-3.0 -1.5 0.0 1.5 3.0

(EII,4,6014) Atom and bond con



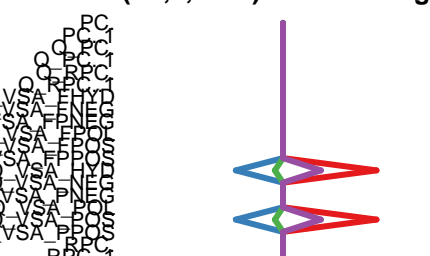
-3.0 -1.5 0.0 1.5 3.0

(EII,4,6014) Atom and bond cou



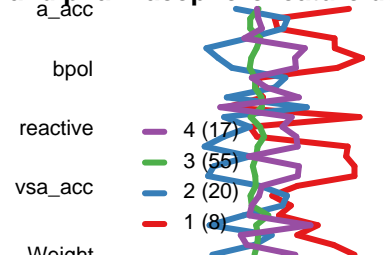
-3.0 -1.5 0.0 1.5 3.0

(EII,4,6014) Partial charge

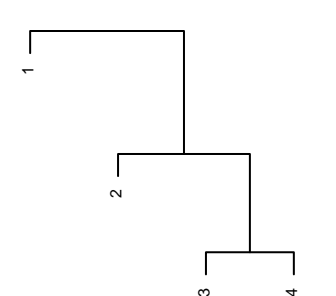
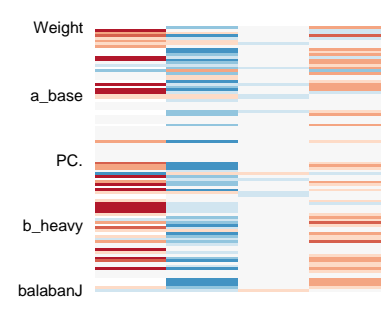


-3.0 -1.5 0.0 1.5 3.0

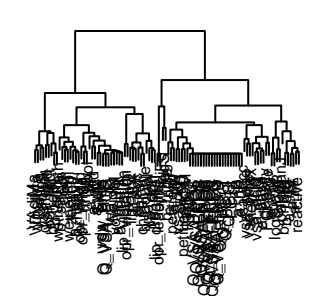
(EII,4,6014) Physical proper and pharmacophore feature des



-3.0 -1.5 0.0 1.5 3.0



dist(avg\_p)  
hclust (\*, "complete")



dist(t(avg\_p))  
hclust (\*, "complete")