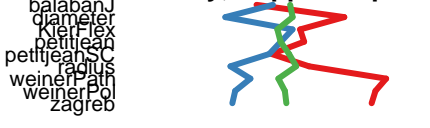
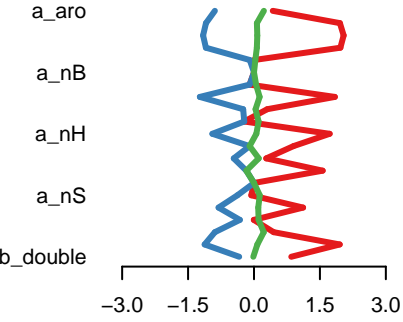


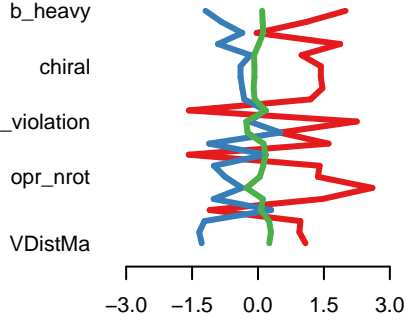
(EII,3,6013) Adjacency, distance matrix, connectivity, Kappa shape in



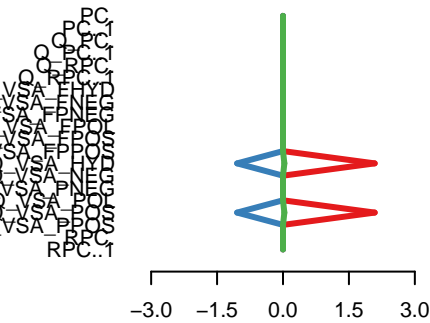
(EII,3,6013) Atom and bond con



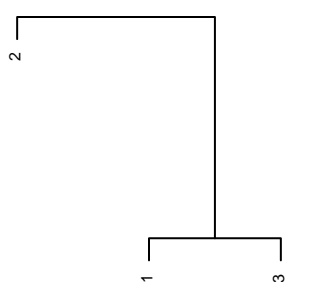
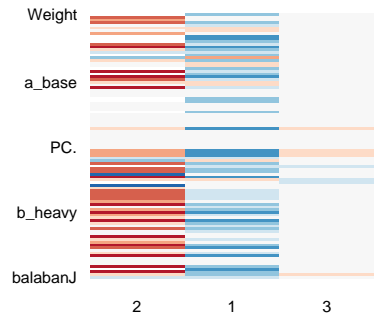
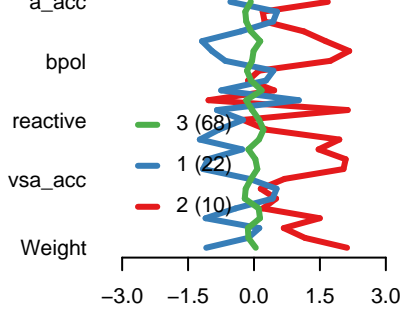
(EII,3,6013) Atom and bond con



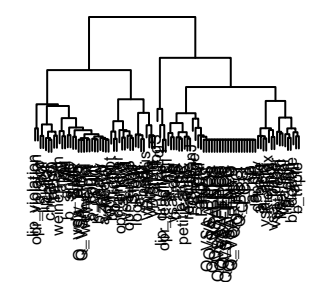
(EII,3,6013) Partial charge



(EII,3,6013) Physical proper and pharmacophore feature des



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



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