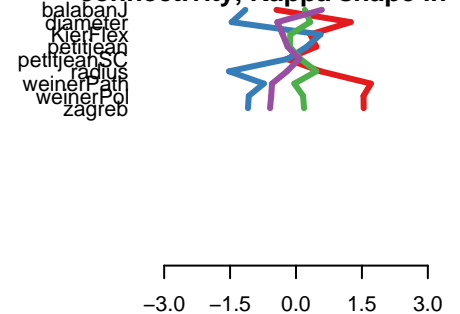
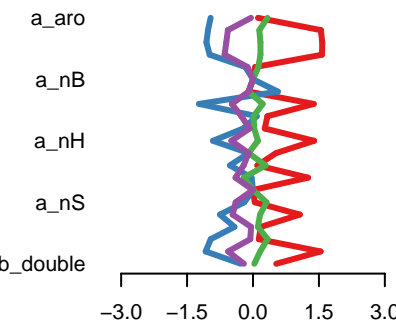


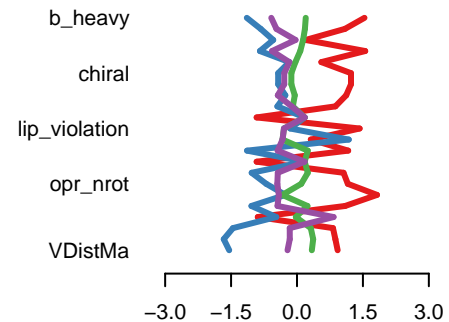
(VII,4,6014) Adjacency, distance matrix, connectivity, Kappa shape in



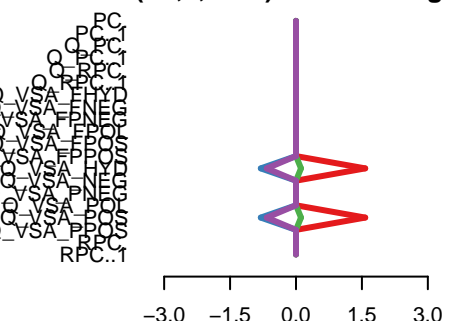
(VII,4,6014) Atom and bond con



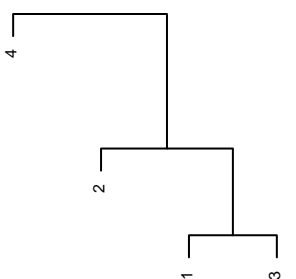
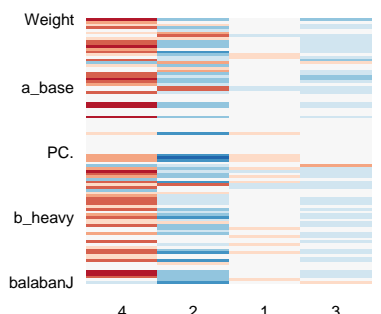
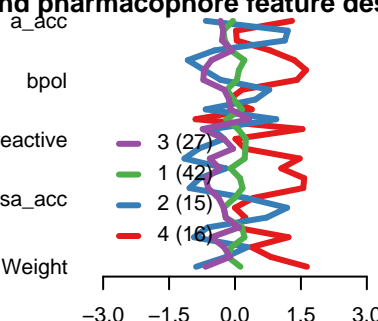
(VII,4,6014) Atom and bond cou



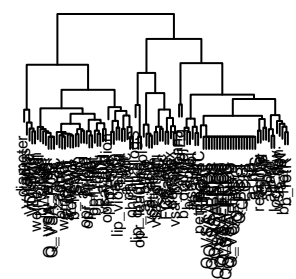
(VII,4,6014) Partial charge



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



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



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