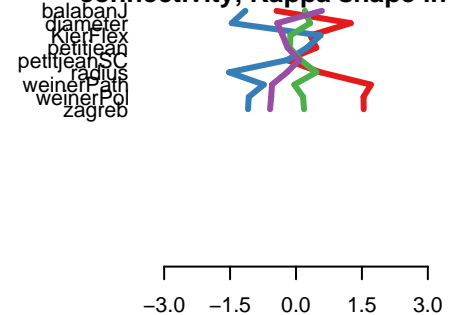
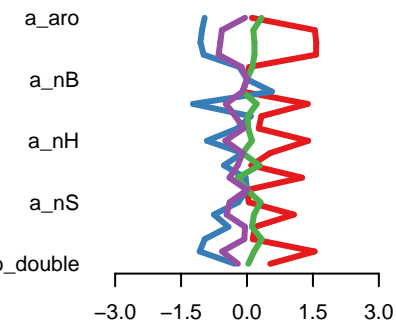


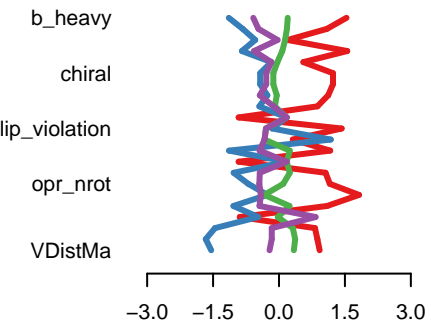
(VII,4,6014) Adjacency, distance matrix, connectivity, Kappa shape in



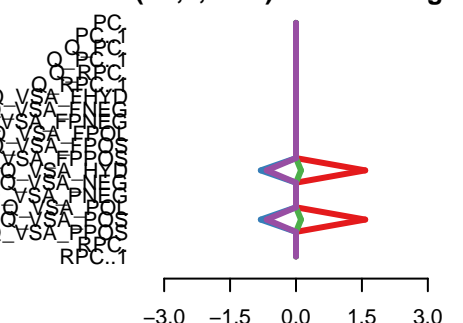
(VII,4,6014) Atom and bond counts



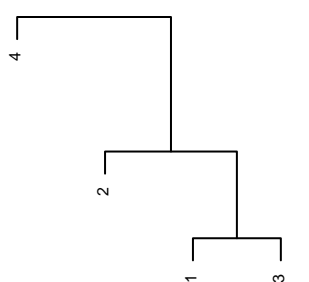
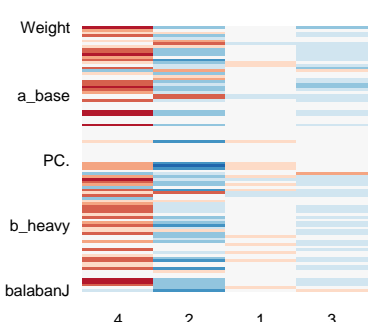
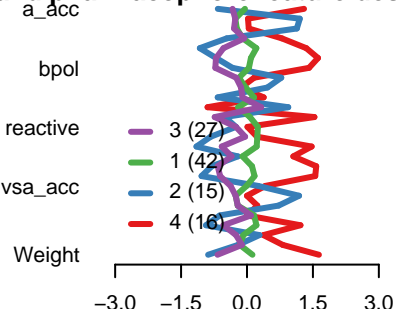
(VII,4,6014) Atom and bond counts



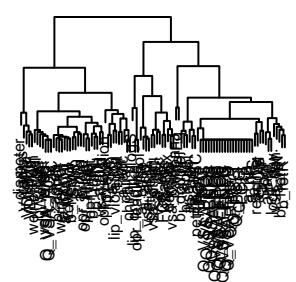
(VII,4,6014) Partial charge



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



dist(avg_p)
hclust (*, "complete")



dist(t(avg_p))
hclust (*, "complete")