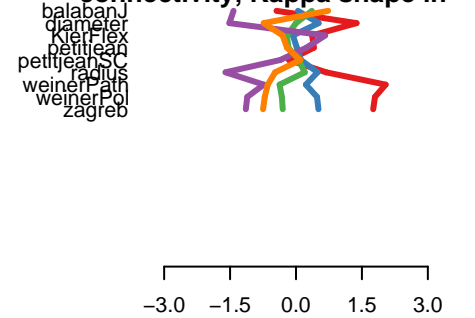
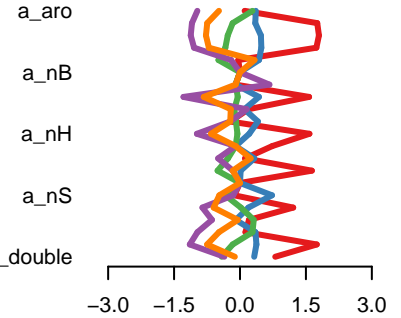


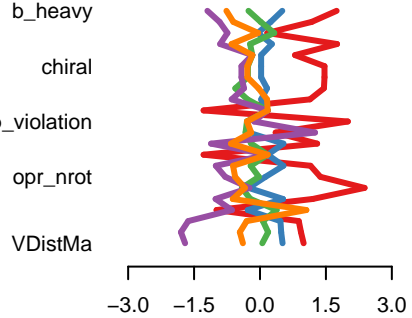
(VII,5,6014) Adjacency, distance matrix, connectivity, Kappa shape in



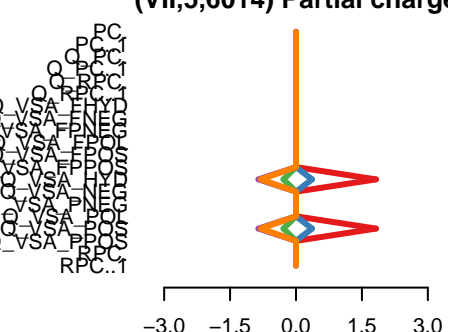
(VII,5,6014) Atom and bond counts



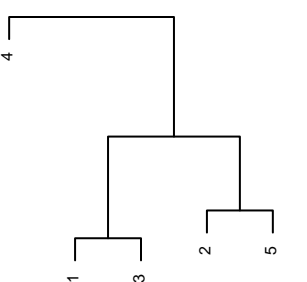
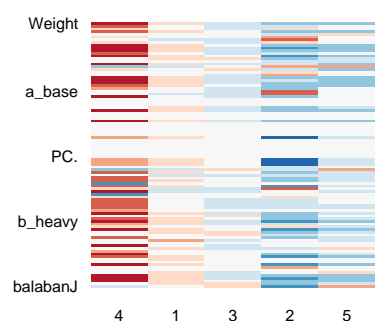
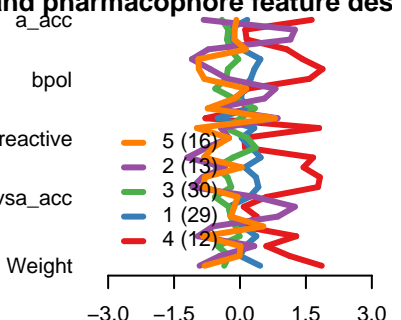
(VII,5,6014) Atom and bond counts



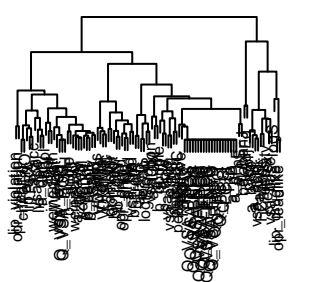
(VII,5,6014) Partial charge



(VII,5,6014) Physical proper and pharmacophore feature descriptors



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



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