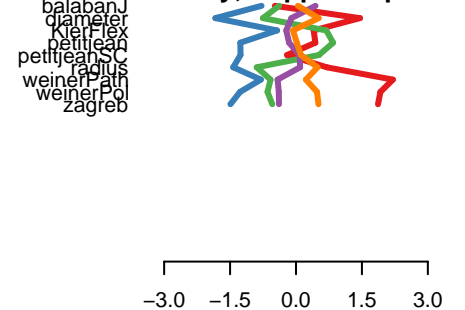
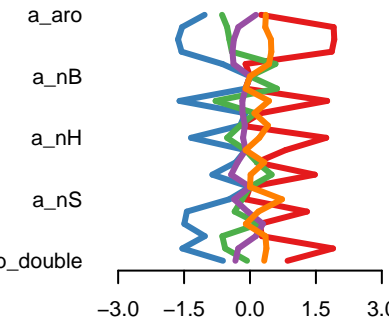


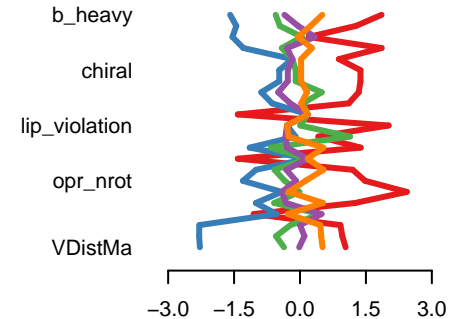
(VII,5,6013) Adjacency, distance matrix, connectivity, Kappa shape in



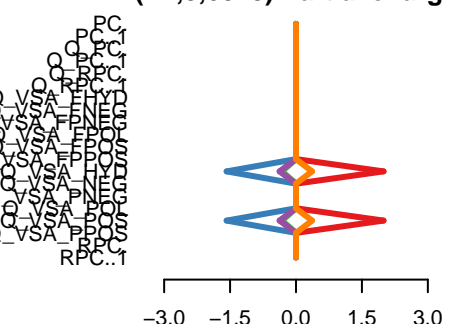
(VII,5,6013) Atom and bond co



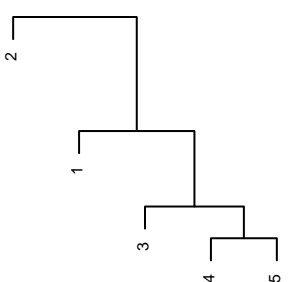
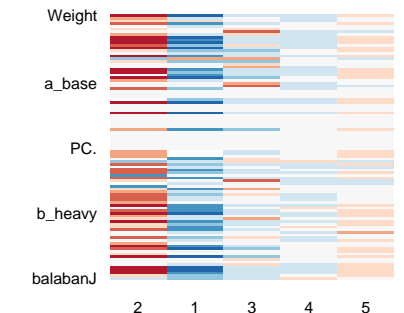
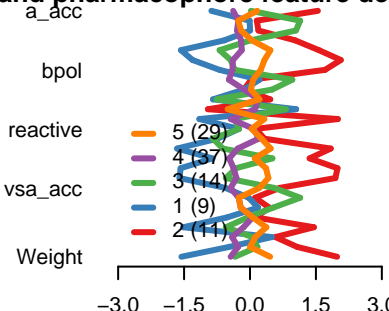
(VII,5,6013) Atom and bond co



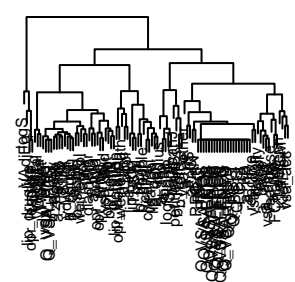
(VII,5,6013) Partial charge



(VII,5,6013) Physical proper and pharmacophore feature des



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



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