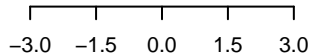
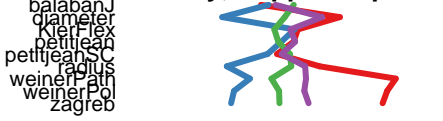
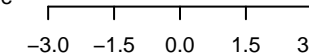
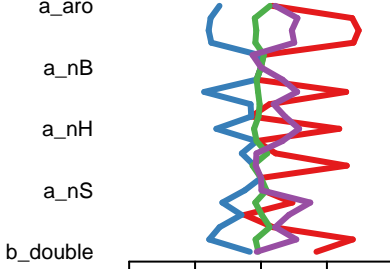


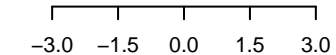
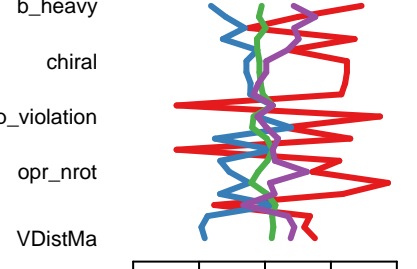
(EII,4,6014) Adjacency, distance matrix, connectivity, Kappa shape in



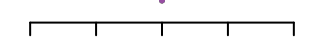
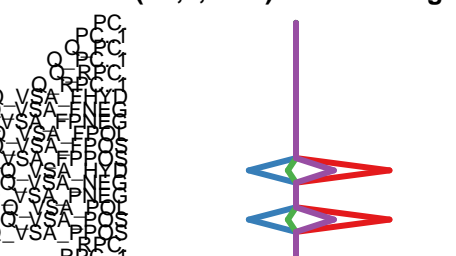
(EII,4,6014) Atom and bond con



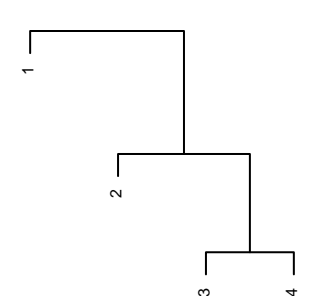
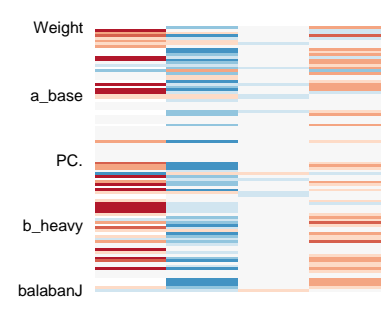
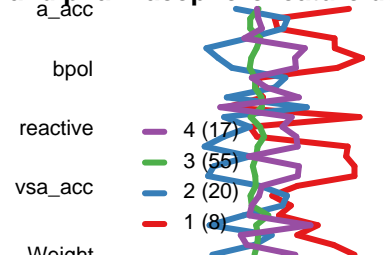
(EII,4,6014) Atom and bond cou



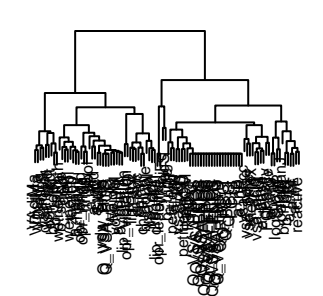
(EII,4,6014) Partial charge



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



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



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