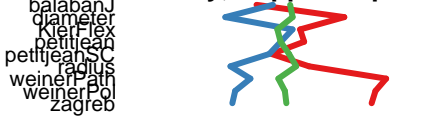
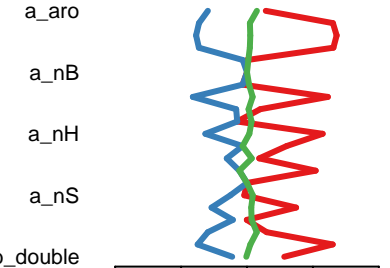


(EII,3,6013) Adjacency, distance matrix, connectivity, Kappa shape in



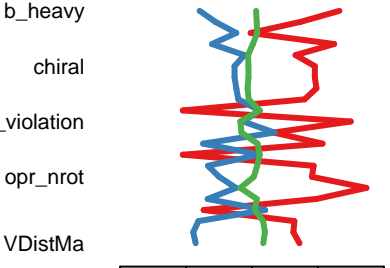
-3.0 -1.5 0.0 1.5 3.0

(EII,3,6013) Atom and bond con



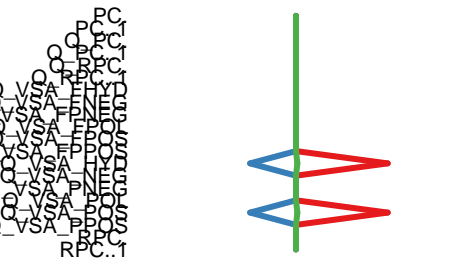
-3.0 -1.5 0.0 1.5 3.0

(EII,3,6013) Atom and bond con



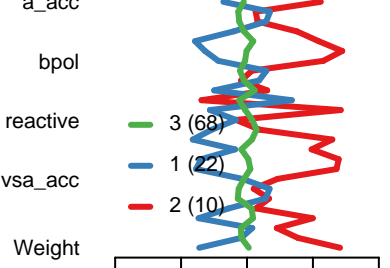
-3.0 -1.5 0.0 1.5 3.0

(EII,3,6013) Partial charge

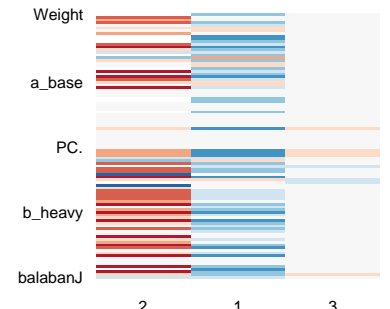


-3.0 -1.5 0.0 1.5 3.0

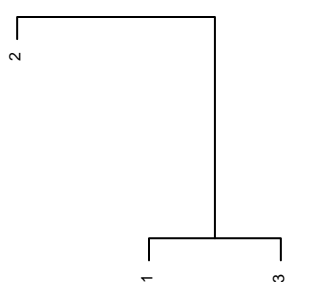
(EII,3,6013) Physical proper and pharmacophore feature des



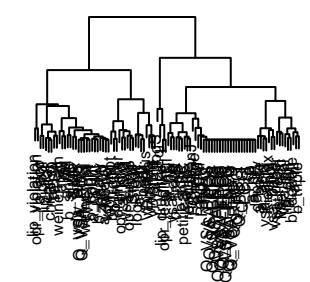
-3.0 -1.5 0.0 1.5 3.0



2 1 3



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



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