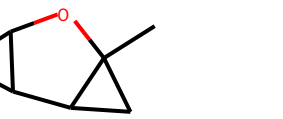
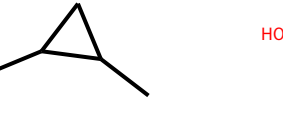


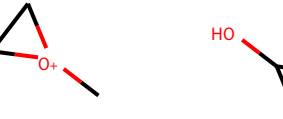
1.00 CC12CC1C1(CC1)O2



0.25 CC12CC1C1CC1O2



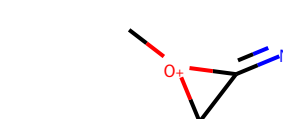
0.11 CC1CC1C1CC1O



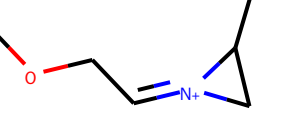
0.11 C[O+]1CC1C1CC1O



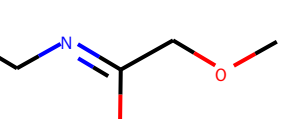
0.11 C[O+]1CC1=C1CC1O



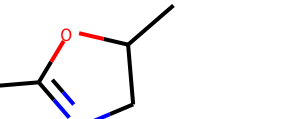
0.11 CC1CC1=C1C[O+]1C



0.11 CC1C[N+]1=C1C[O+]1C



0.07 COCC=[N+]1CC1C



0.02 CCCN=C(O)COC



0.09 COCC1=NCC(C)O1