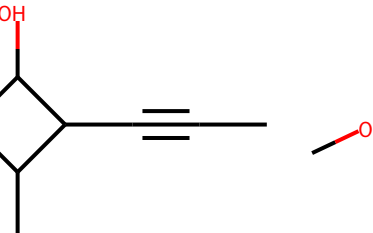
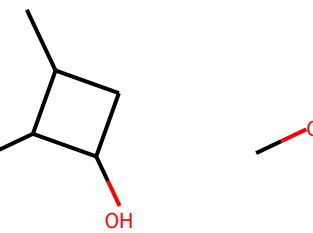


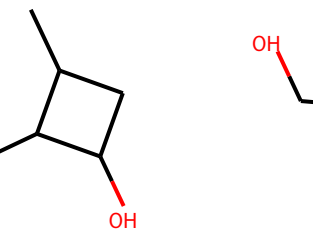
1.00 CC#CC1(CC)CC1O



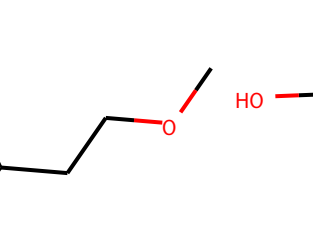
0.27 CC#CC1C(C)CC1O



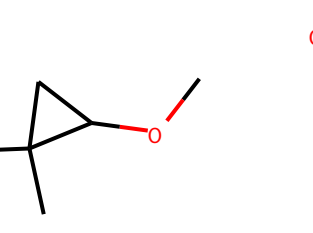
0.15 C[O+]=CC1C(C)CC1O



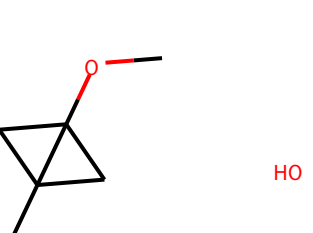
0.17 COCC1C(C)CC1O



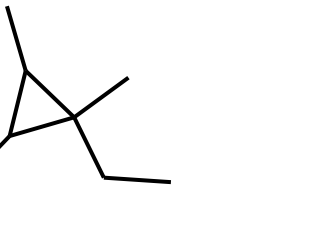
0.09 COCCC(C)CCO



0.21 COC1CC1(C)CCO



0.17 COC12CC1(C)C2CO



0.21 CCC1(C)C(C)C1CO