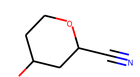
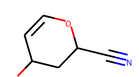
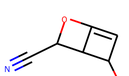
1.00 N#CC12CC(O)C1CO2



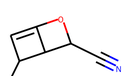
0.25 N#CC1CC(O)CCO1



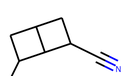
0.22 N#CC1CC(O)C=CO1



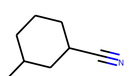
0.19 N#CC1OC2=CC(O)C21



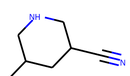
0.14 CC1C=C2OC(C#N)C21



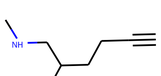
0.16 CC1CC2CC(C#N)C12



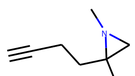
0.16 CC1CCCC(C#N)C1



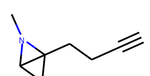
0.16 CC1CNCC(C#N)C1



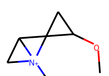
0.02 C#CCCC(C)CNC



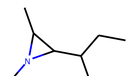
0.07 C#CCCC1(C)CN1C



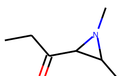
0.11 C#CCCC12CC1N2C



0.14 COC1CC12C1C[N+]12C



0.04 CCC(C)C1C(C)N1C



0.04 CCC(=O)C1C(C)N1C