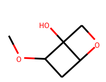
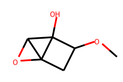
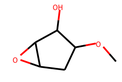


1.00 COC1CC20CC12O



0.35 COC1CC23OC2C13O



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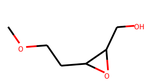
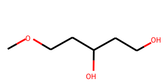
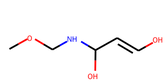


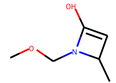
0.16 COCCC1OC1CO



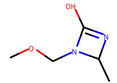
0.09 COCCC(O)CCO



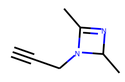
0.09 COCNC(O)C=CO



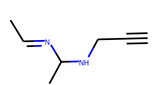
0.11 COCN1C(O)=CC1C



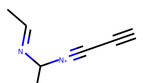
0.11 COCN1C(O)=NC1C



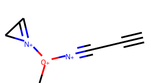
0.04 C#CCN1C(C)=NC1C



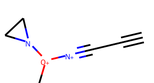
0.02 C#CCNC(C)N=CC



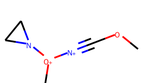
0.02 C#CC#[N+][C@H](C)N=CC



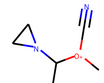
0.04 C#CC#[N+][O+](C)[N+]1=CC1



0.07 C#CC#[N+][O+](C)N1CC1



0.11 COC#[N+][O+](C)N1CC1



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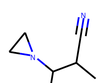

0.07 CC(C#N)C(C)N1CC1