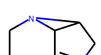
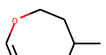


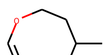
1.00 C1CN2C3CN=C(O1)C32

0.38 C1CC2C3CN=C(O1)C23

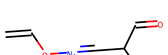
0.17 C1=C2CC(CC1)CCO2



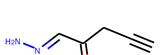
0.13 CC1CCC=COCC1



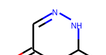
0.10 CC1C=CC=COCC1



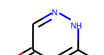
0.00 C=CO[N+]#CC(C)C=O



0.00 C#CCC(=O)C=NN



0.09 CC1CC(=O)C=NN1



0.04 Cc1cc(=O)cn[nH]1