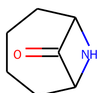
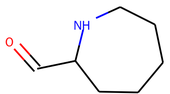
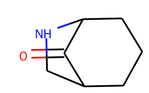
1.00 O=C1C2CC=CCC1N2



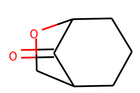
0.37 O=C1C2CCCCC1N2



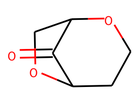
0.11 O=CC1CCCCN1



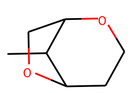
0.24 O=C1C2CCCC1NC2



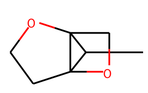
0.18 O=C1C2CCCC1OC2



0.18 O=C1C2CCOC1CO2



0.08 CC1C2CCOC1CO2



0.06 CC1C23CCOC12CO3