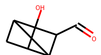
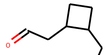
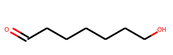


1.00 O=CC1C2C3CC2C13O

0.26 O=CC1C(O)C2CCC12



0.17 O=CCC1CCC1CO



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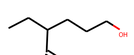

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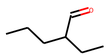

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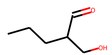

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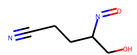
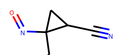
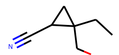


0.04 N#CCCC(CO)N=O



0.12 N#CC1CC1(CO)N=O



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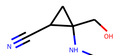

0.09 CNC1(CO)CC1C#N