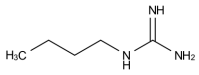
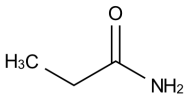
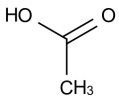
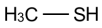
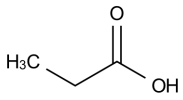
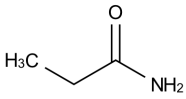
# Chemical Library Test

| 2D Molecule | SMILES |
|---|---|
|  | C |
|  | CCCCNC(N)=N |
|  | CCC(N)=O |
|  | CC(O)=O |
|  | CS |
|  | CCC(O)=O |
|  | CCC(N)=O |

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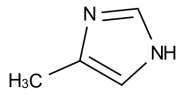
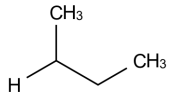
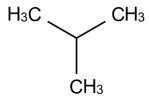
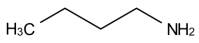
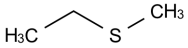
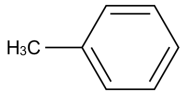
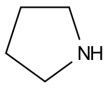
|---|---|
| H | [H] |
|   | CC1=CNC=N1 |
|   | C(CC)([H])C |
|   | CC(C)C |
|   | CCCCN |
|   | CCSC |
|   | CC1=CC=CC=C1 |
|   | C2CCCN2 |

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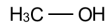
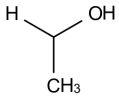
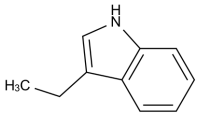
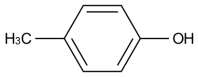
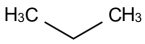
|---|---|
|  | CO |
|  | C(C)([H])O |
|  | CCC1=CNC2=C1C=CC=C2 |
|  | CC1=CC=C(O)C=C1 |
|  | C(C)C |