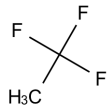
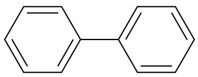
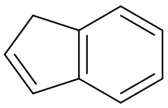
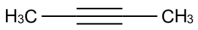
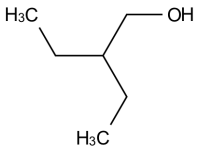
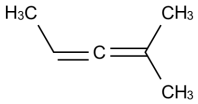
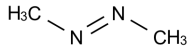
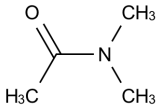
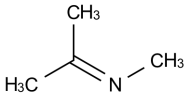
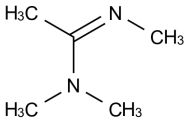
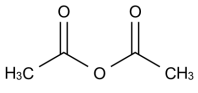
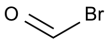
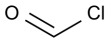
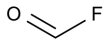
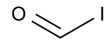
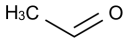
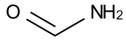
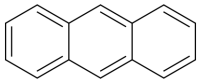
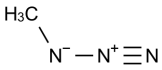
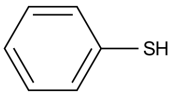
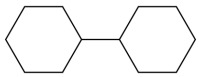
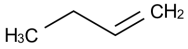
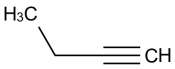
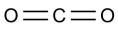
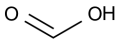
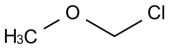
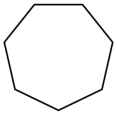
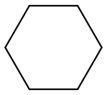
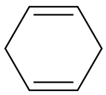
Functional Groups Global Chem

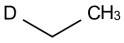
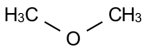
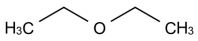
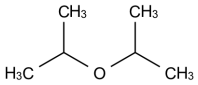
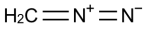
2D Molecule	SMILES
	CC(F)(F)F
	C1(C2=CC=CC=C2)=CC=CC=C1
	C1(CC=C2)=C2C=CC=C1
	[NH]1CCCC1
	CC#CC
	CCC(CC)CO
	CC=C=C(C)C

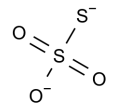
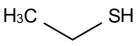
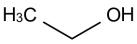
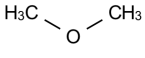
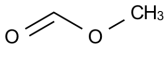
	C/N=N/C
	CC(N(C)C)=O
	C/C(C)=N/C
	C/C(N(C)C)=N/C
	CC(=O)OC(=O)C
	C(=O)Br
	C(=O)Cl
	C(=O)F

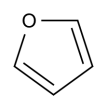
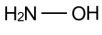
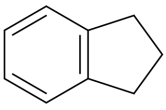
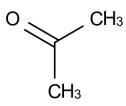
	C(=O)I
	CC=O
	C(=O)N
	*N
	C12=CC=CC=C1C=C3C(C=CC=C3)=C2
	C([N-][N+]#N)
	C1=CC=CC=C1
	C1=CC=C(C=C1)S

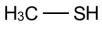
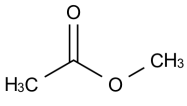
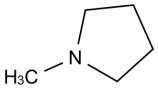
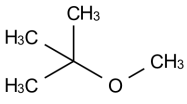
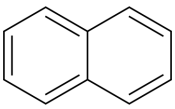
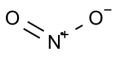
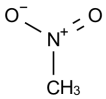
	C1CCCCC1C1CCCCC1
HBr	Br
	CCC=C
	CCC#C
	O=C=O
	C(=O)O
HCl	Cl
	COCCl

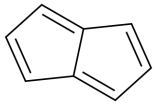
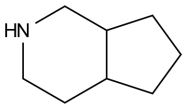
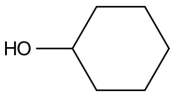
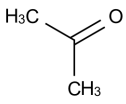
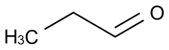
	C1=CC=C1
	C1CCC1
	C1CCCCCC1
	C1CCCCC1
	C1=CCCC=C1
	C1=CCC=CC1
	C=1CCCCC=1
	C1CCCC1

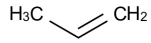
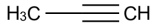
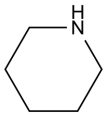
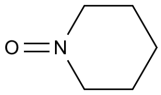
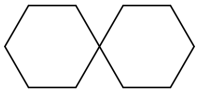
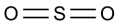
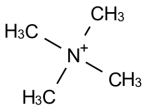
	C1=CCC=C1
	C1CC1
	C1=CC1
	[2H][CH2]C
	COC
	CCOCC
	CC(C)OC(C)C
Failed Rendering	C&1&1&1&1
	C=[N+]=[N-]

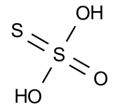
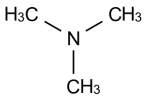
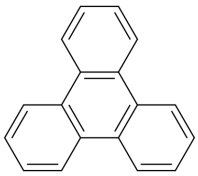
 NH_4^+ NH_4^+	[NH4+].[NH4+].[O-]S(=O)(=O)[S-]
NH_3	N
H_3C-CH_3	CC
	CCS
	CCO
$H_2C=CH_2$	C=C
	COC
	C(=O)OC

HF	F
	C=O
	C1OC=CC=1
Failed Rendering	C&1&1&1
	C#N
	[OH-]
	NO
	C1=CC=CC(CCC2)=C12
	CC(=O)C

	C
	CS
	CC(OC)=O
	CN1CCCC1
	CC(C)(C)OC
	C12=CC=CC=C1C=CC=C2
	[N+](=O)[O-]
	C[N+]([O-])=O

	C12=CC=CC1=CC=C2
	N1CC2CCCC2CC1
	OC1CCCCC1
	C=1(C=CC=CC1)
Failed Rendering	c1ccccc1C&1&1
H_2O	O
NH_3	N
	CC(C)=O
	CCC=O

	CC=C
	CC#C
	N1CCCCC1
	O=N1CCCCC1
	NC
	C12(CCCCC1)CCCCC2
	S(=O)(=O)
	C[N+](C)(C)C

H_2S	S
	OS(=O)(=S)O
	CN(C)C
	C1(C=CC=C2)=C2C(C=CC=C3)=C3C4=C1C=CC=C4