Functional Groups Global Chem

2D Molecule	SMILES
F F F H₃C	CC(F)(F)F
	C1(C2=CC=CC)=CC=C1
	C1(CC=C2)=C2C=CC=C1
NH	[NH]1CCCC1
H ₃ C ————— CH ₃	CC#CC
H ₃ C OH	CCC(CC)CO
H ₃ C CH ₃ CCH ₃	CC=C=C(C)C
H₃C N CH₃	C/N=N/C

O CH ₃ N H ₃ C CH ₃	CC(N(C)C)=O
CH ₃ CH ₃	C/C(C)=N/C
H_3C N CH_3 H_3C N CH_3	C/C(N(C)C)=N/C
H₃C O CH₃	CC(=O)OC(=O)C
O Br	C(=O)Br
O CI	C(=O)CI
0 F	C(=O)F
0	C(=O)I

H₃CO	CC=O
O NH ₂	C(=O)N
A — NH ₂	*N
	C12=CC=CC=C1C=C3C(C=CC=C3)=C2
H_3C $N^N^+\equiv N$	C([N-][N+]#N)
	C1=CC=CC1
SH	C1=CC=C(C=C1)S
	C1CCCCC1C1CCCCC1

HBr	Br
H ₃ C CH ₂	CCC=C
H ₃ CCH	CCC#C
o=c=0	O=C=O
О	C(=O)O
HCI	CI
H₃C CI	COCCI
	C1=CC=C1

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

	C1CC1
	C1=CC1
D CH₃	[2H][CH2]C
H ₃ C CH ₃	coc
H₃C CH₃	ccocc
H₃C CH₃ CH₃ CH₃	CC(C)OC(C)C
Failed Rendering	C&1&1&1&1
$H_2C = N^+ = N^-$	C=[N+]=[N-]
O S S O O O O O O O O O O O O O O O O O	[NH4+].[NH4+].[O-]S(=O)(=O)[S-]

NH ₃	N
H₃C — CH₃	СС
H₃C SH	ccs
H₃C OH	cco
$H_2C = CH_2$	C=C
H₃C CH₃	сос
0	C(=O)OC
HF	F

H₂C == O	C=O
	C1OC=CC=1
Failed Rendering	C&1&1&1
HC≡N	C#N
HO ⁻	[OH-]
H₂N — OH	NO
	C1=CC=CC(CCC2)=C12
O CH ₃	CC(=O)C
CH₄	С

H₃C — SH	CS
H ₃ C CH ₃	CC(OC)=O
H ₃ C N	CN1CCCC1
H₃C CH₃ CH₃ CH₃	CC(C)(C)OC
	C12=CC=CC=C1C=CC=C2
O	[N+](=O)[O-]
O O O O O O O O O O O O O O O O O O O	C[N+]([O-])=O
	C12=CC=CC1=CC=C2

HN	N1CC2CCC2CC1
но —	OC1CCCCC1
	C=1(C=CC=CC1)
Failed Rendering	c1ccccc1C&1&1
H ₂ O	0
NH₃	N
H ₃ C O CH ₃	CC(C)=O
H ₃ C O	CCC=O
H₃C CH₂	CC=C

H ₃ C — — CH	CC#C
✓—K	N1CCCCC1
o=N	O=N1CCCCC1
H₂N — CH₃	NC
	C12(CCCCC1)CCCCC2
o=s=o	S(=O)(=O)
H ₃ C / CH ₃ / CH ₃	C[N+](C)(C)C
H ₂ S	S

OH S / S / HO	OS(=O)(=S)O
H ₃ C CH ₃	CN(C)C
	C1(C=CC=C2)=C2C(C=CC=C3)=C3C4=C1C=CC=C4