Input coordinates for 2.

Standard orientation:

Center Number	Atomic Number	Atomic Type	Coordinates (Angstroms) X	Y	Z
1	6	0	3.068038	1.130984	0.015002
2	6	0	3.338171	-0.200198	-0.003290
3	6	0	1.967053	-2.315391	-0.034397
4	7	0	1.687857	1.257760	0.012022
5	7	0	2.115285	-0.858011	-0.015753
6	17	0	-0.643794	0.092922	2.197039
7	6	0	0.962597	2.528670	0.038886
8	6	0	1.101089	0.038212	-0.005203
9	23	0	-0.999539	-0.374611	-0.006359
10	8	0	-1.045273	-1.930201	-0.032758
11	17	0	-3.051078	0.469249	0.017784
12	17	0	-0.662174	0.167531	-2.194861
13	1	0	3.724183	1.986461	0.027207
14	1	0	4.275961	-0.731878	-0.010920
15	1	0	1.404724	-2.622940	-0.916700
16	1	0	1.442695	-2.650438	0.861601
17	1	0	2.963158	-2.757964	-0.062770
18	1	0	0.393387	2.610732	0.967393
19	1	0	0.287872	2.584056	-0.817715
20	1	0	1.685695	3.342802	-0.018997