

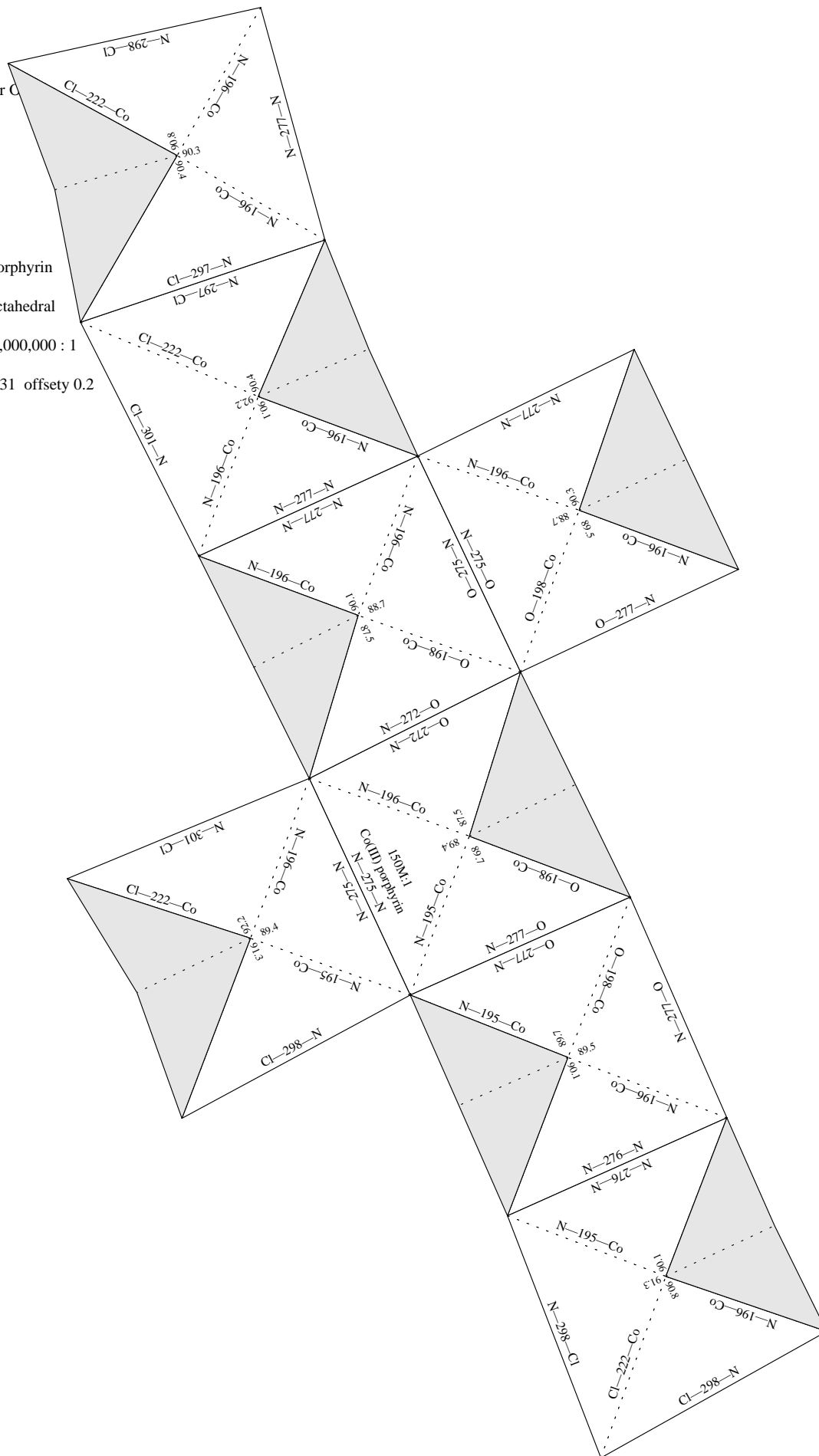
Molecular Origami

given information

ElementNames	[(Co) (Cl) (N) (N) (N) (N) (O)]	
distance	195.010	Co ¹ -N ¹
distance	195.545	Co ¹ -N ²
distance	195.780	Co ¹ -N ³
distance	195.844	Co ¹ -N ⁴
distance	197.841	Co ¹ -O ¹
distance	221.620	Co ¹ -Cl ¹
angle	87.484	O ¹ -Co ¹ -N ⁴
	272.2	O ¹ -N ⁴
angle	88.699	O ¹ -Co ¹ -N ³
	275.2	O ¹ -N ³
angle	89.406	N ⁴ -Co ¹ -N ¹
	274.9	N ⁴ -N ¹
angle	89.543	O ¹ -Co ¹ -N ²
	277.1	O ¹ -N ²
angle	89.676	O ¹ -Co ¹ -N ¹
	277.	O ¹ -N ¹
angle	90.090	N ² -Co ¹ -N ¹
	276.4	N ² -N ¹
angle	90.118	N ⁴ -Co ¹ -N ³
	277.2	N ⁴ -N ³
angle	90.302	N ³ -Co ¹ -N ²
	277.4	N ³ -N ²
angle	90.360	N ³ -Co ¹ -Cl ¹
	296.6	N ³ -Cl ¹
angle	90.828	N ² -Co ¹ -Cl ¹
	297.7	N ² -Cl ¹
angle	91.263	N ¹ -Co ¹ -Cl ¹
	298.4	N ¹ -Cl ¹
angle	92.153	N ⁴ -Co ¹ -Cl ¹
	301.2	N ⁴ -Cl ¹
angle	176.987	N ⁴ -Co ¹ -N ²
	391.3	N ⁴ -N ²
angle	178.325	N ³ -Co ¹ -N ¹
	390.7	N ³ -N ¹
angle	178.990	O ¹ -Co ¹ -Cl ¹
	419.4	O ¹ -Cl ¹
dopage	T	
AutoAlign	F	

structure type: XABCDEF

View -1



actual size: 400 709