Molecular Origami of Co(diethylenetriamine)Cl3facial

given information

ElementNames	[(Co) (Cl) (Cl) (Cl) (N) (N) (N)]	
distance	196.736	Co^1-N^1
distance	197.630	Co^1-N^3
distance	200.876	Co^1-N^2
distance	224.312	Co^1-Cl^1
distance	227.748	Co^1-Cl^3
distance	230.912	Co^1-Cl^2
angle	84.335	N^3-Co^1-Cl^2
	288.7	N^3-Cl^2
angle	85.710	N^1-Co^1-Cl^2
	291.9	N^1-Cl^2
angle	87.479	N^2-Co^1-Cl^3
	297.	N^2-Cl^3
angle	88.792	N^3-Co^1-Cl^3
	298.4	N^3-Cl^3
angle	89.075	N^1-Co^1-Cl^3
	298.5	N^1-Cl^3
angle	90.013	N^3-Co^1-Cl^1
	299.	N^3-Cl^1
angle	90.340	N^2-Co^1-Cl^1
	302.	N^2-Cl^1
angle	91.005	Cl^3-Co^1-Cl^2
	327.2	Cl^3-Cl^2
angle	91.204	Cl^2-Co^1-Cl^1
	325.3	Cl^2-Cl^1
angle	92.507	N^1-Co^1-Cl^1
	304.8	N^1-Cl^1
angle	93.123	N^2-Co^1-N^1
	288.7	N^2-N^1
angle	96.767	N^3-Co^1-N^2
	297.9	N^3-N^2
angle	169.779	N^3-Co^1-N^1
	392.8	N^3-N^1
angle	177.374	Cl^3-Co^1-Cl^1
	451.9	Cl^3-Cl^1
angle	178.104	N^2-Co^1-Cl^2
	431.7	N^2-Cl^2
dopage	T	
AutoAlign	F	

structure type: XABCDEF

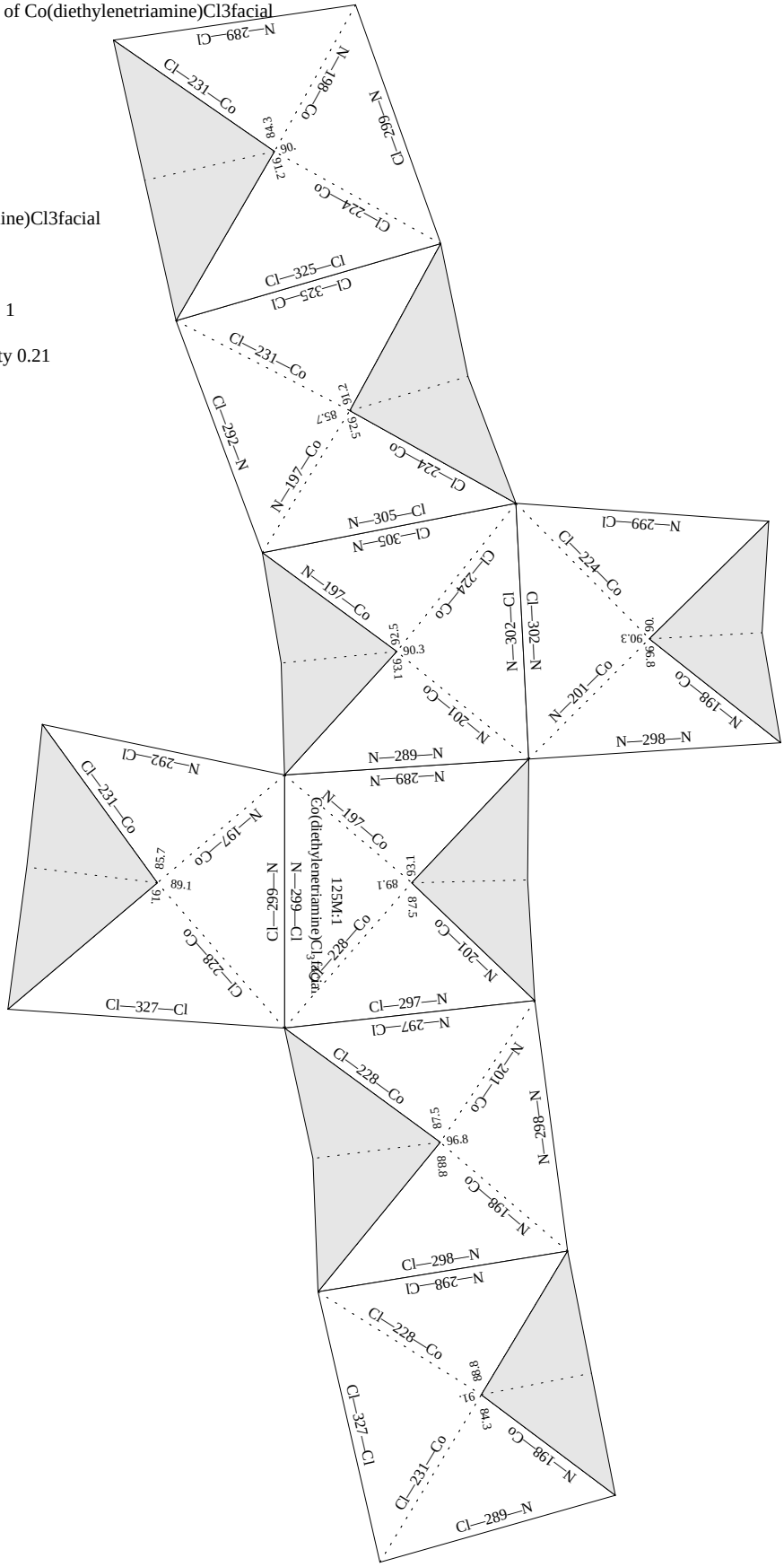
Molecular Origami of Co(diethylenetriamine)Cl3facial

!Co1
Cl1
Cl2
Cl3
N1
N2
N3
Co(diethylenetriamine)Cl3facial

special AX6?

scale 125,000,000 : 1
units: pm
offsetx -0.05 offsety 0.21

View -1



Current: (centerx 4.25) (centery 5.21) (scale 125)
%%BoundingBox: 127 81 470 752 actual: 137 91 460 742 center: 299 417 actual size: 323 651
Better: (centerx 4.35) (centery 4.92) (scale 125)
%%BoundingBox: 131 76 474 746 actual: 141 86 464 736 center: 302 411 actual size: 323 651