Molecular Origami of SbPhCl2

given information

ElementNames	[(Sb) (C) (Cl) (Cl)]	
dotted	F	
distance	215.051	Sb^1-C^1
distance	237.558	Sb^1-Cl^2
distance	241.097	Sb^1-Cl^1
angle	92.603	Cl^1-Sb^1-C^1
	330.3	Cl^1-C^1
angle	94.349	Cl^2-Sb^1-Cl^1
	351.1	Cl^2-Cl^1
angle	95.291	Cl^2-Sb^1-C^1
	334.8	Cl^2-C^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

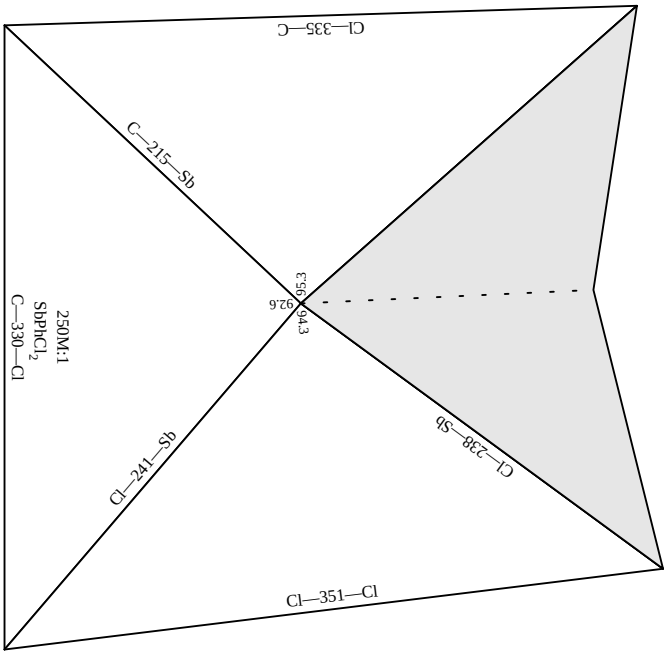
Molecular Origami of SbPhCl2

!Sb1
C1
Cl1
Cl2
SbPhCl2

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx 0.0 offsety -0.81

View -1



Current: (centerx 4.30) (centery 4.19) (scale 250)
%%BoundingBox: 183 292 450 553 actual: 193 302 440 543 center: 316 422 actual size: 247 241
Better: (centerx 4.16) (centery 3.82) (scale 250)
%%BoundingBox: 173 207 439 468 actual: 183 217 429 458 center: 306 338 actual size: 247 241