Molecular Origami of Sb(C(CH3)3)Cl2

given information

ElementNames	[(Sb) (C) (Cl) (Cl)]	
dotted	F	
distance	219.021	Sb^1-C^1
distance	237.304	Sb^1-Cl^2
distance	238.776	Sb^1-Cl^1
angle	95.024	Cl^2-Sb^1-C^1
	336.7	Cl^2-C^1
angle	95.152	Cl^1-Sb^1-C^1
	338.2	Cl^1-C^1
angle	95.398	Cl^2-Sb^1-Cl^1
	352.1	Cl^2-Cl^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

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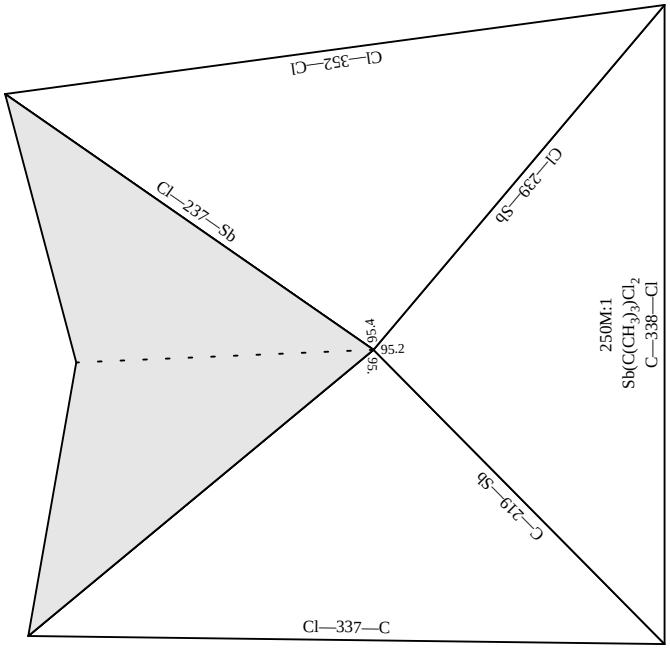
!Sb1
C1
Cl1
Cl2
Sb(C(CH3)3)Cl2

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx -0.05 offsety 2.62

View -1



Current: (centerx 4.25) (centery 7.62) (scale 250)
%%BoundingBox: 169 299 436 559 actual: 179 309 426 549 center: 302 429 actual size: 247 240
Better: (centerx 4.30) (centery 7.16) (scale 250)
%%BoundingBox: 169 455 436 714 actual: 179 465 426 704 center: 302 585 actual size: 247 240