Molecular Origami of Sb((CH3)3CCH2)3

given information

ElementNames	[(Sb) (C) (C) (C)]	
dotted	F	
distance	218.307	Sb^{1}-C^{7}
distance	218.310	Sb^{1}-C^{1}
distance	218.311	Sb^{1}-C^{6}
angle	93.502	C^{7}-Sb^{1}-C^{1}
	318.	C^{7}-C^{1}
angle	93.502	C^{7}-Sb^{1}-C^{6}
	318.	C^{7}-C^{6}
angle	93.504	C^{6}-Sb^{1}-C^{1}
	318.	C^{6}-C^{1}
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

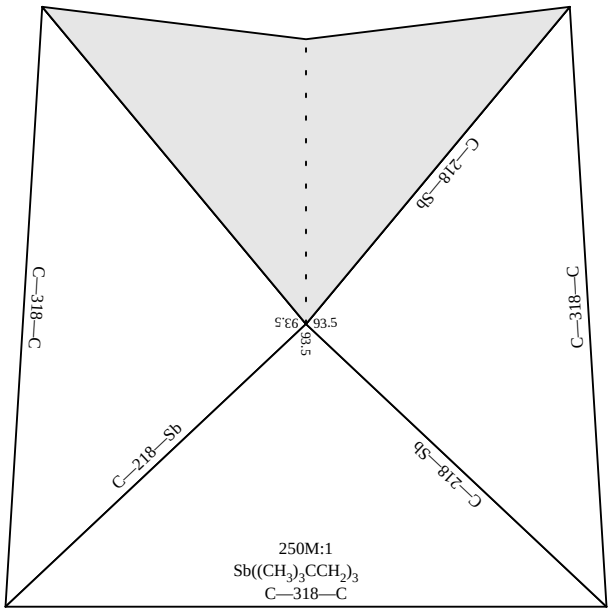
Molecular Origami of Sb((CH3)3CCH2)3

!Sb1
C1
C6
C7
Sb((CH3)3CCH2)3

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx 1.51 offsety 0.9

View -1



Current: (centerx 5.81) (centery 5.90) (scale 250)
%%BoundingBox: 183 302 428 547 actual: 193 312 418 537 center: 306 425 actual size: 225 225
Better: (centerx 5.82) (centery 5.50) (scale 250)
%%BoundingBox: 292 338 537 583 actual: 302 348 527 573 center: 415 461 actual size: 225 225