Molecular Origami of Sb(pCH3Ph)Br2

given information

ElementNames	[(Sb) (Br) (Br) (C)]	
dotted	F	
distance	215.575	Sb^1-C^1
distance	251.581	Sb^1-Br^1
distance	255.051	Sb^1-Br^2
angle	92.746	C^1-Sb^1-Br^2
	341.7	C^1-Br^2
angle	95.181	C^1-Sb^1-Br^1
	345.8	C^1-Br^1
angle	95.729	Br^2-Sb^1-Br^1
	375.7	Br^2-Br^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

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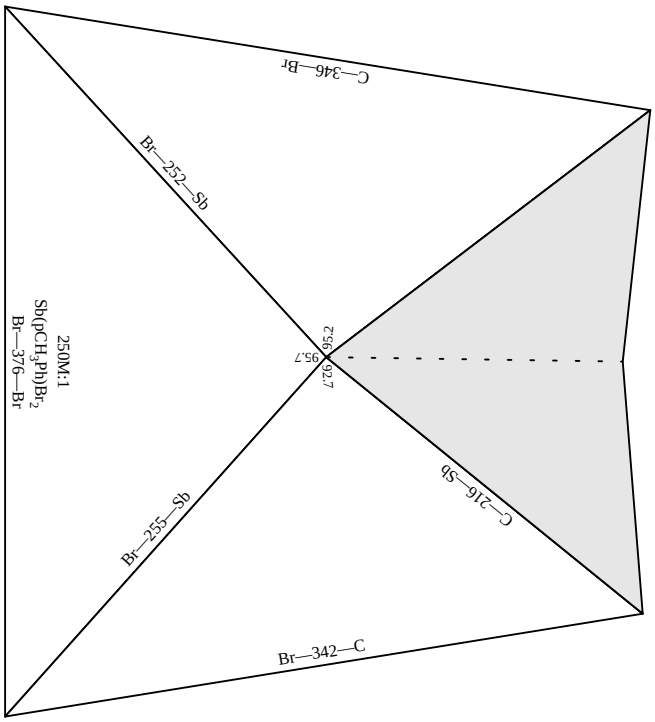
!Sb1
Br1
Br2
C1
Sb(pCH3Ph)Br2

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx -0.19 offsety -0.93

View -1



Current: (centerx 4.11) (centery 4.07) (scale 250)
%%BoundingBox: 153 283 415 569 actual: 163 293 405 559 center: 284 426 actual size: 242 266
Better: (centerx 4.42) (centery 3.65) (scale 250)
%%BoundingBox: 161 186 423 472 actual: 171 196 413 462 center: 292 329 actual size: 242 266