Molecular Origami of SbPh2Br

given information

ElementNames	[(Sb) (Br) (C) (C)]	
dotted	F	
distance	214.250	Sb^1-C^1
distance	214.980	Sb^1-C^7
distance	255.226	Sb^1-Br^1
angle	93.476	C^7-Sb^1-Br^1
	343.5	C^7-Br^1
angle	95.284	C^1-Sb^1-Br^1
	348.	C^1-Br^1
angle	98.515	C^7-Sb^1-C^1
	325.2	C^7-C^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

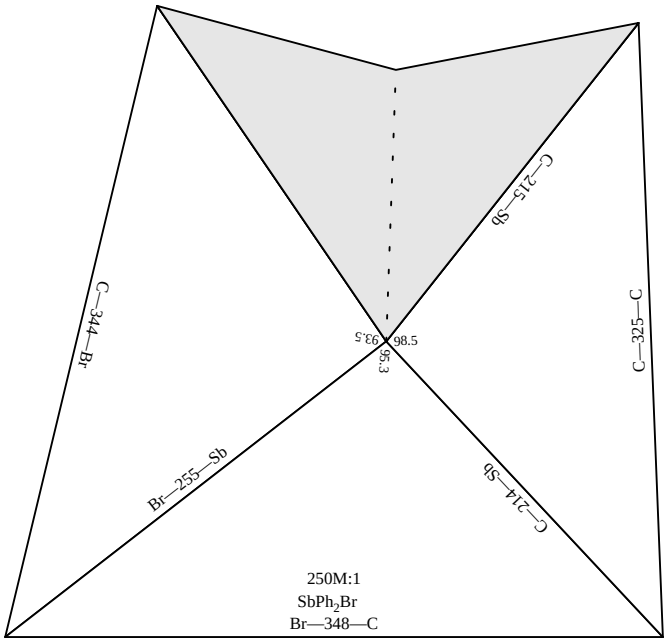
Molecular Origami of SbPh2Br

!Sb1
Br1
C1
C7
SbPh2Br

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx 1.59 offsety 0.9

View -1



Current: (centerx 5.89) (centery 5.90) (scale 250)
%%BoundingBox: 167 291 434 548 actual: 177 301 424 538 center: 301 420 actual size: 247 237
Better: (centerx 5.96) (centery 5.57) (scale 250)
%%BoundingBox: 287 332 554 589 actual: 297 342 544 579 center: 420 461 actual size: 247 237