Molecular Origami of SbPhBr2

given information

ElementNames	[(Sb) (Br) (Br) (C)]	
dotted	F	
distance	215.864	Sb^1-C^1
distance	252.575	Sb^1-Br^2
distance	256.287	Sb^1-Br^1
angle	92.959	C^1-Sb^1-Br^1
	343.5	C^1-Br^1
angle	95.562	Br^2-Sb^1-Br^1
	376.9	Br^2-Br^1
angle	96.100	C^1-Sb^1-Br^2
	349.3	C^1-Br^2
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

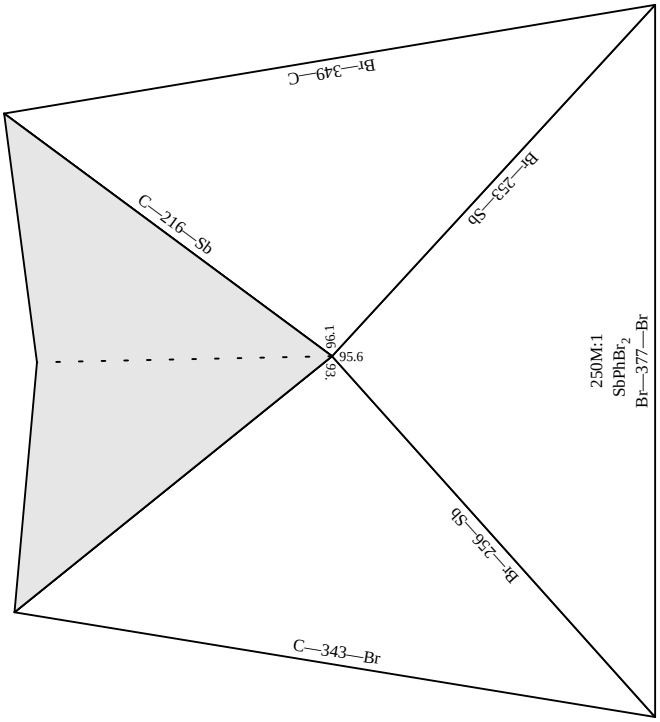
Molecular Origami of SbPhBr2

!Sb1
Br1
Br2
C1
SbPhBr2

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx 0.07 offsety 2.73

View -1



Current: (centerx 4.37) (centery 7.73) (scale 250)
%%BoundingBox: 194 279 458 567 actual: 204 289 448 557 center: 326 423 actual size: 244 267
Better: (centerx 4.09) (centery 7.35) (scale 250)
%%BoundingBox: 179 449 443 736 actual: 189 459 433 726 center: 311 593 actual size: 244 267