Molecular Origami of SbBr3

given information

ElementNames	[(Sb) (Br) (Br) (Br)]	
dotted	F	
distance	250.073	Sb^{1}-Br^{3}
distance	250.574	Sb^{1}-Br^{1}
distance	252.134	Sb^{1}-Br^{2}
angle	93.403	Br^{3}-Sb^{1}-Br^{1}
	364.4	Br^{3}-Br^{1}
angle	95.588	Br^{3}-Sb^{1}-Br^{2}
	372.	Br^{3}-Br^{2}
angle	96.483	Br^{2}-Sb^{1}-Br^{1}
	375.	Br^{2}-Br^{1}
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

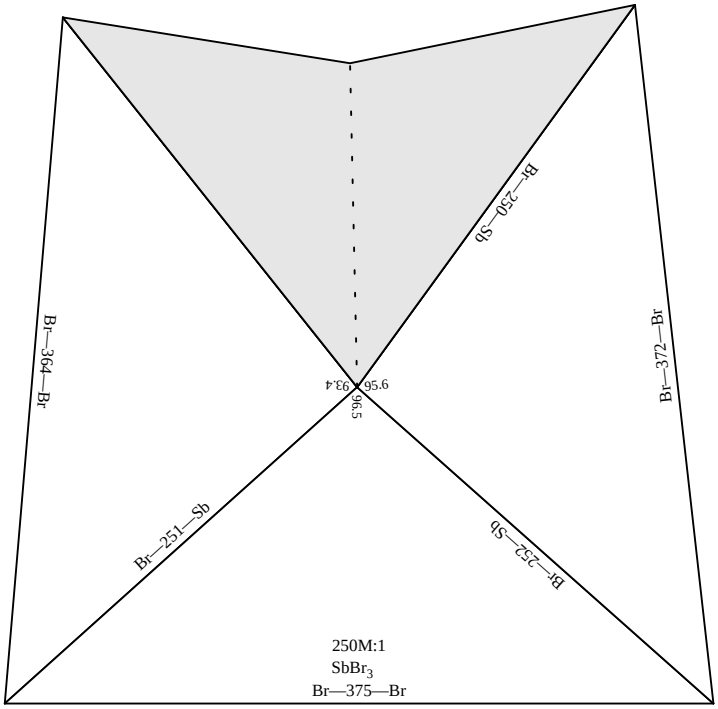
Molecular Origami of SbBr3

!Sb1
Br1
Br2
Br3
SbBr3

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx 1.77 offsety 0.9

View -1



Current: (centerx 6.07) (centery 5.90) (scale 250)
%%BoundingBox: 161 282 447 564 actual: 171 292 437 554 center: 304 423 actual size: 266 262
Better: (centerx 6.10) (centery 5.53) (scale 250)
%%BoundingBox: 291 320 576 602 actual: 301 330 566 592 center: 433 461 actual size: 266 262