Molecular Origami of SbBr3

given information

ElementNames	[(Sb) (Br) (Br) (Br)]	
dotted	F	
distance	248.982	Sb^{1}-Br^{3}
distance	250.252	Sb^{1}-Br^{1}
distance	250.524	Sb^{1}-Br^{2}
angle	93.925	Br^{2}-Sb^{1}-Br^{1}
	366.	Br^{2}-Br^{1}
angle	94.020	Br^{3}-Sb^{1}-Br^{1}
	365.2	Br^{3}-Br^{1}
angle	97.263	Br^{3}-Sb^{1}-Br^{2}
	374.9	Br^{3}-Br^{2}
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

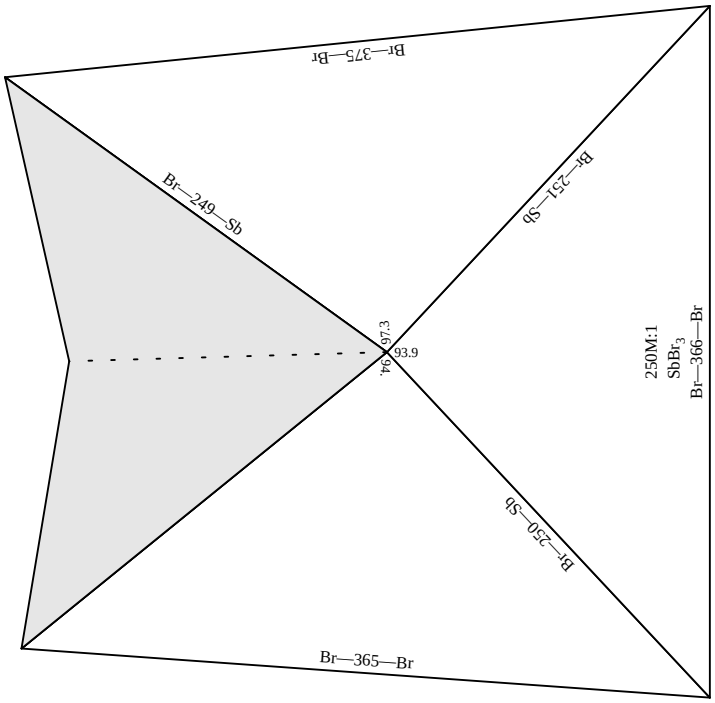
Molecular Origami of SbBr3

!Sb1
Br1
Br2
Br3
SbBr3

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx -0.05 offsety 2.74

View -1



Current: (centerx 4.25) (centery 7.74) (scale 250)
%%BoundingBox: 161 288 446 567 actual: 171 298 436 557 center: 304 428 actual size: 264 259
Better: (centerx 4.28) (centery 7.30) (scale 250)
%%BoundingBox: 160 454 445 733 actual: 170 464 435 723 center: 302 593 actual size: 264 259