Molecular Origami of SbBr3

given information

ElementNames	[(Sb) (Br) (Br) (Br)]	
dotted	F	
distance	248.993	Sb^{1}-Br^{1}
distance	248.999	Sb^{1}-Br^{2}
distance	249.004	Sb^{1}-Br^{3}
angle	98.000	Br^{3}-Sb^{1}-Br^{1}
	375.8	Br^{3}-Br^{1}
angle	98.001	Br^{3}-Sb^{1}-Br^{2}
	375.9	Br^{3}-Br^{2}
angle	98.002	Br^{2}-Sb^{1}-Br^{1}
	375.8	Br^{2}-Br^{1}
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

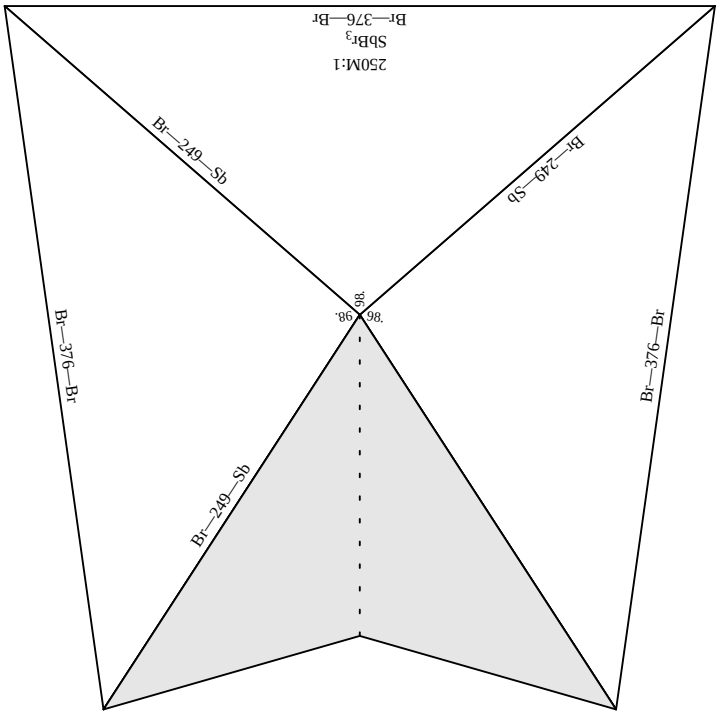
Molecular Origami of SbBr3

!Sb1
Br1
Br2
Br3
SbBr3

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx -1.88 offsety 0.9

View -1



Current: (centerx 2.42) (centery 5.90) (scale 250)
%%BoundingBox: 164 284 451 568 actual: 174 294 441 558 center: 307 426 actual size: 266 264
Better: (centerx 2.40) (centery 5.48) (scale 250)
%%BoundingBox: 27 319 314 603 actual: 37 329 304 593 center: 171 461 actual size: 266 264