Molecular Origami of $SbPh_3$

given information

ElementNames	[(Sb) (C) (C) (C)]	
dotted	F	
distance	214.287	Sb^{1}-C^{1}
distance	215.071	Sb^{1}-C^{13}
distance	215.326	Sb^{1}-C^{7}
angle	95.657	C^{13}-Sb^{1}-C^{1}
	318.2	C^{13}-C^{1}
angle	96.005	C^{7}-Sb^{1}-C^{13}
	319.9	C^{7}-C^{13}
angle	98.022	C^{7}-Sb^{1}-C^{1}
	324.3	C^{7}-C^{1}
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

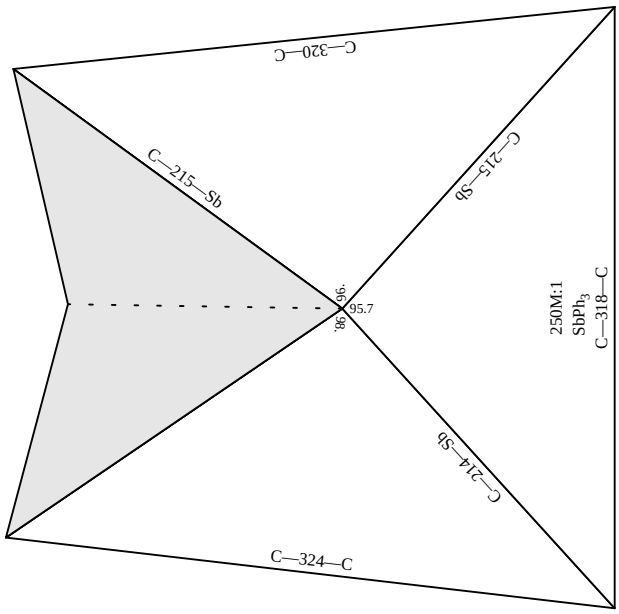
Molecular Origami of SbPh3

!Sb1
C1
C13
C7
SbPh3

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx -0.05 offsety 2.49

View -1



Current: (centerx 4.25) (centery 7.49) (scale 250)
%%BoundingBox: 180 304 429 549 actual: 190 314 419 539 center: 305 427 actual size: 228 226
Better: (centerx 4.27) (centery 7.07) (scale 250)
%%BoundingBox: 178 453 427 698 actual: 188 463 417 688 center: 302 575 actual size: 228 226