Molecular Origami of SbPhI2

given information

ElementNames	[(Sb) (C) (I) (I)]	
dotted	F	
distance	213.853	Sb^1-C^1
distance	273.809	Sb^1-I^2
distance	275.332	Sb^1-I^1
angle	94.498	I^1-Sb^1-C^1
	361.6	I^1-C^1
angle	97.705	I^2-Sb^1-I^1
	413.5	I^2-I^1
angle	97.796	I^2-Sb^1-C^1
	369.6	I^2-C^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

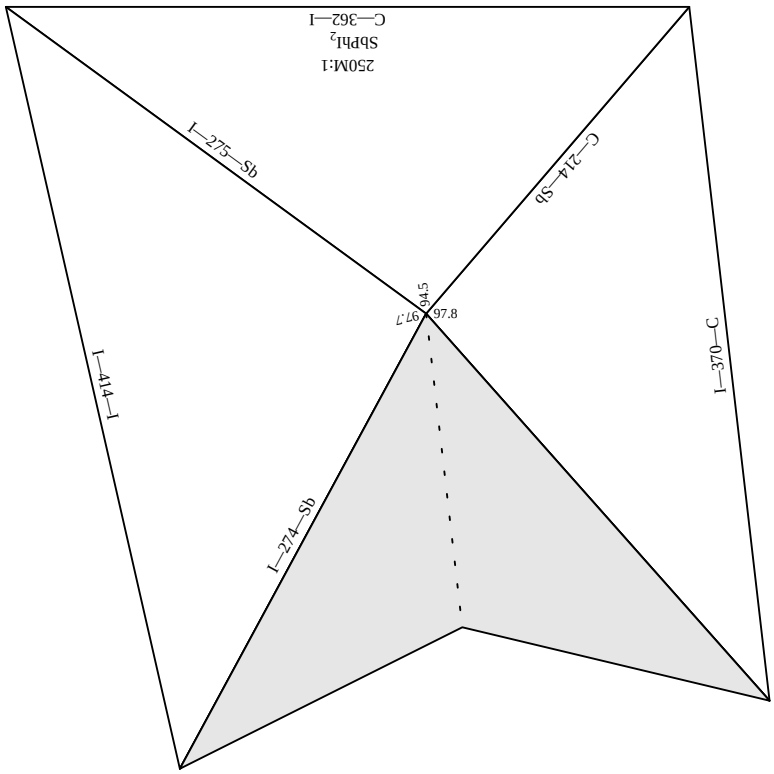
Molecular Origami of SbPhI2

!Sb1
C1
I1
I2
SbPhI2

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx -2.03 offsety 0.7

View -1



Current: (centerx 2.27) (centery 5.70) (scale 250)
%%BoundingBox: 153 243 460 549 actual: 163 253 450 539 center: 307 396 actual size: 286 286
Better: (centerx 2.26) (centery 5.70) (scale 250)
%%BoundingBox: 7 294 313 599 actual: 17 304 303 589 center: 160 446 actual size: 286 286