Molecular Origami of As(Ph)3

given information

ElementNames	[(As) (C) (C) (C)]	
dotted	F	
distance	193.510	As^{1}-C^{7}
distance	195.776	As^{1}-C^{13}
distance	196.100	As^{1}-C^{1}
angle	99.216	C^{13}-As^{1}-C^{1}
	298.5	C^{13}-C^{1}
angle	99.571	C^{7}-As^{1}-C^{13}
	297.3	C^{7}-C^{13}
angle	100.376	C^{7}-As^{1}-C^{1}
	299.3	C^{7}-C^{1}
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

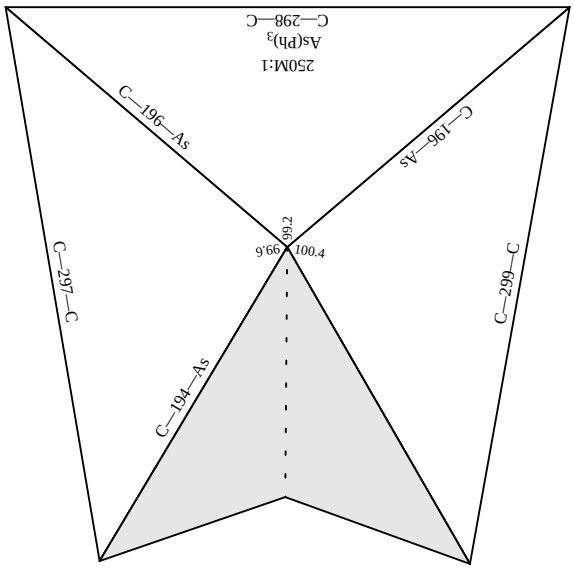
Molecular Origami of As(Ph)3

!As1
C1
C13
C7
As(Ph)3

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx -1.5 offsety 0.9

View -1



Current: (centerx 2.80) (centery 5.90) (scale 250)
%%BoundingBox: 192 312 423 541 actual: 202 322 413 531 center: 307 426 actual size: 212 209
Better: (centerx 2.78) (centery 5.48) (scale 250)
%%BoundingBox: 82 346 314 575 actual: 92 356 304 565 center: 198 461 actual size: 212 209