Molecular Origami of PBu3

given information

ElementNames	[(P) (C) (C) (C)]	
dotted	F	
distance	184.016	P^1-C^5
distance	184.375	P^1-C^1
distance	184.606	P^1-C^9
angle	100.503	C^5-P^1-C^1
	283.2	C^5-C^1
angle	102.182	C^9-P^1-C^1
	287.1	C^9-C^1
angle	102.688	C^9-P^1-C^5
	287.9	C^9-C^5
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

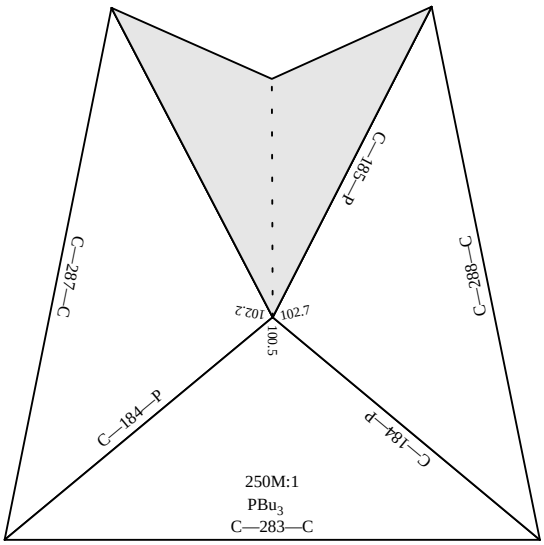
Molecular Origami of PBu3

!P1
C1
C5
C9
PBu3

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx 1.34 offsety 0.9

View -1



Current: (centerx 5.64) (centery 5.90) (scale 250)
%%BoundingBox: 195 314 416 534 actual: 205 324 406 524 center: 306 424 actual size: 201 200
Better: (centerx 5.64) (centery 5.51) (scale 250)
%%BoundingBox: 292 351 513 571 actual: 302 361 503 561 center: 402 461 actual size: 201 200