Molecular Origami of P(CH(CH3)2)3

given information

ElementNames	[(P) (C) (C) (C)]	
dotted	F	
distance	185.636	P^1-C^7
distance	185.986	P^1-C^1
distance	187.043	P^1-C^4
angle	101.984	C^7-P^1-C^4
	289.6	C^7-C^4
angle	102.907	C^7-P^1-C^1
	290.6	C^7-C^1
angle	104.703	C^4-P^1-C^1
	295.4	C^4-C^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

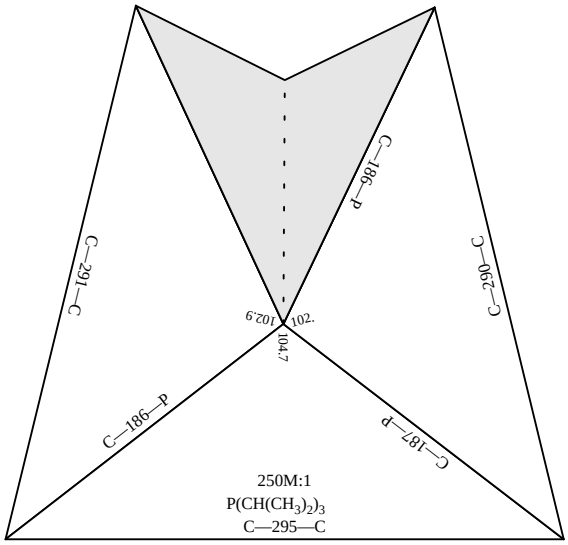
Molecular Origami of P(CH(CH3)2)3

!P1
C1
C4
C7
P(CH(CH3)2)3

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx 1.36 offsety 0.9

View -1



Current: (centerx 5.66) (centery 5.90) (scale 250)
%%BoundingBox: 188 310 418 530 actual: 198 320 408 520 center: 303 420 actual size: 209 200
Better: (centerx 5.70) (centery 5.56) (scale 250)
%%BoundingBox: 289 351 519 571 actual: 299 361 509 561 center: 404 461 actual size: 209 200