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

given information

ElementNames	[(P) (C) (C) (C)]	
dotted	F	
distance	183.181	P^1-C^1
distance	184.993	P^1-C^7
distance	186.189	P^1-C^{10}
angle	100.982	C^{10}-P^1-C^1
	285.	C^{10}-C^1
angle	102.817	C^7-P^1-C^{10}
	290.1	C^7-C^{10}
angle	104.345	C^7-P^1-C^1
	290.8	C^7-C^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

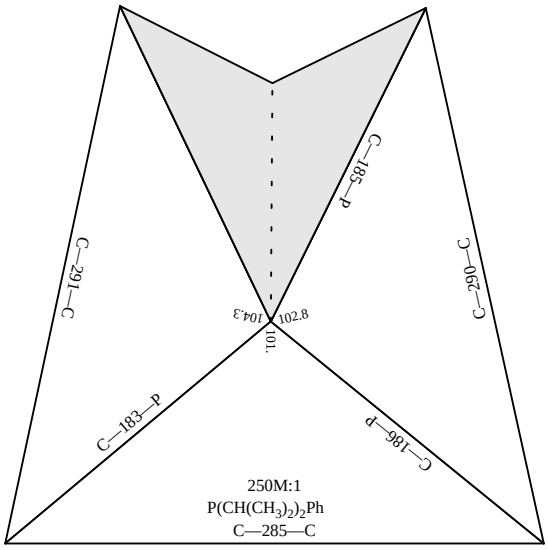
Molecular Origami of P(CH(CH3)2)2Ph

!P1
C1
C10
C7
P(CH(CH3)2)2Ph

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx 1.35 offsety 0.9

View -1



Current: (centerx 5.65) (centery 5.90) (scale 250)
%%BoundingBox: 195 314 417 535 actual: 205 324 407 525 center: 306 425 actual size: 202 202
Better: (centerx 5.65) (centery 5.50) (scale 250)
%%BoundingBox: 292 350 514 572 actual: 302 360 504 562 center: 403 461 actual size: 202 202