

# Molecular Origami of P(C6H11)3

given information

ElementNames	[ (P) (C) (C) (C) ]	
dotted	F	
distance	186.590	P <sup>1</sup> -C <sup>7</sup>
distance	186.637	P <sup>1</sup> -C <sup>1</sup>
distance	187.186	P <sup>1</sup> -C <sup>13</sup>
angle	103.150	C <sup>13</sup> -P <sup>1</sup> -C <sup>1</sup>
	292.9	C <sup>13</sup> -C <sup>1</sup>
angle	103.228	C <sup>7</sup> -P <sup>1</sup> -C <sup>13</sup>
	293.	C <sup>7</sup> -C <sup>13</sup>
angle	105.058	C <sup>7</sup> -P <sup>1</sup> -C <sup>1</sup>
	296.2	C <sup>7</sup> -C <sup>1</sup>
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

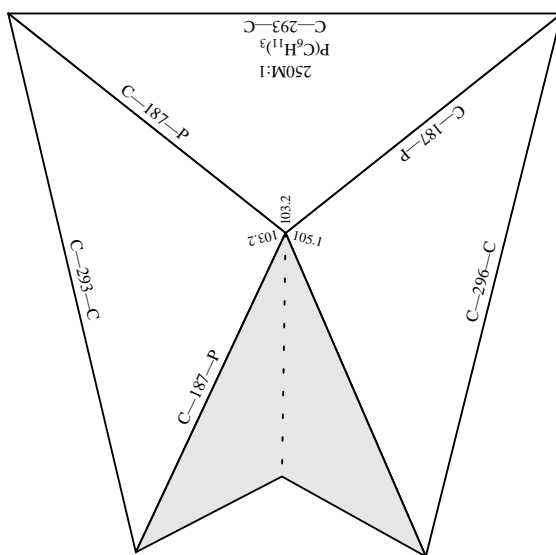
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!P1  
C1  
C13  
C7  
P(C6H11)3

trigonal pyramidal

scale 250,000,000 : 1  
units: pm  
offsetx -1.45 offsety 0.9

View -1



Current: (centerx 2.85) (centery 5.90) (scale 250)

%%BoundingBox: 195 315 423 539 actual: 205 325 413 529

center: 309 427

actual size: 208 204

Better: (centerx 2.81) (centery 5.47) (scale 250)

%%BoundingBox: 88 349 315 573

actual: 98 359 305 563

center: 202 461

actual size: 208 204