

Molecular Origami of P(CCH)<sub>3</sub>

given information

ElementNames	[ (P) (C) (C) (C) ]	
dotted	F	
distance	175.011	P <sup>1</sup> -C <sup>5</sup>
distance	175.011	P <sup>1</sup> -C <sup>1</sup>
distance	176.316	P <sup>1</sup> -C <sup>3</sup>
angle	98.819	C <sup>5</sup> -P <sup>1</sup> -C <sup>1</sup>
	265.8	C <sup>5</sup> -C <sup>1</sup>
angle	101.827	C <sup>5</sup> -P <sup>1</sup> -C <sup>3</sup>
	272.7	C <sup>5</sup> -C <sup>3</sup>
angle	101.827	C <sup>3</sup> -P <sup>1</sup> -C <sup>1</sup>
	272.7	C <sup>3</sup> -C <sup>1</sup>
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

# Molecular Origami of P(CCH)<sub>3</sub>

!P1

C1

C3

C5

P(CCH)<sub>3</sub>

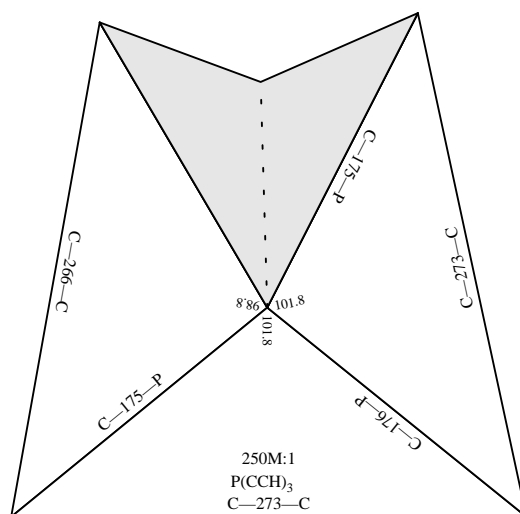
trigonal pyramidal

scale 250,000,000 : 1

units: pm

offsetx 1.26 offsety 0.9

View -1



Current: (centerx 5.56) (centery 5.90) (scale 250)

%%BoundingBox: 197 318 410 527 actual: 207 328 400 517

Better: (centerx 5.59) (centery 5.53) (scale 250)

%%BoundingBox: 290 356 503 565 actual: 300 366 493 555

center: 304 423

actual size: 193 189

center: 397 461

actual size: 193 189