Molecular Origami of P(Ph2)2CH2

given information

ElementNames	[(P) (C) (C) (C)]	
dotted	F	
distance	182.998	P^1-C^2
distance	184.250	P^1-C^8
distance	184.573	P^1-C^1
angle	101.246	C^8-P^1-C^1
	285.1	C^8-C^1
angle	101.711	C^8-P^1-C^2
	284.8	C^8-C^2
angle	103.192	C^2-P^1-C^1
	288.	C^2-C^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

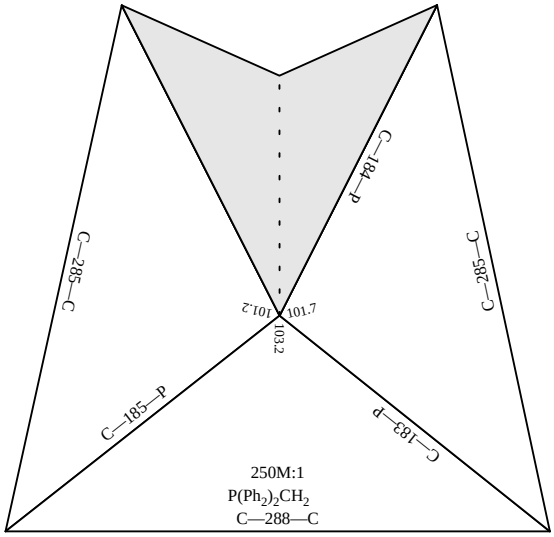
Molecular Origami of P(Ph2)2CH2

!P1
C1
C2
C8
P(Ph2)2CH2

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: 192 313 416 530 actual: 202 323 406 520 center: 304 421 actual size: 204 197
Better: (centerx 5.67) (centery 5.55) (scale 250)
%%BoundingBox: 290 352 515 569 actual: 300 362 505 559 center: 402 461 actual size: 204 197