Molecular Origami of P((CH3O)2Ph)3

given information

ElementNames	[(P) (C) (C) (C)]	
dotted	F	
distance	183.654	P^1-C^1
distance	184.023	P^1-C^9
distance	185.411	P^1-C^{17}
angle	99.255	C^9-P^1-C^1
	280.1	C^9-C^1
angle	106.703	C^{17}-P^1-C^1
	296.1	C^{17}-C^1
angle	107.273	C^9-P^1-C^{17}
	297.5	C^9-C^{17}
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

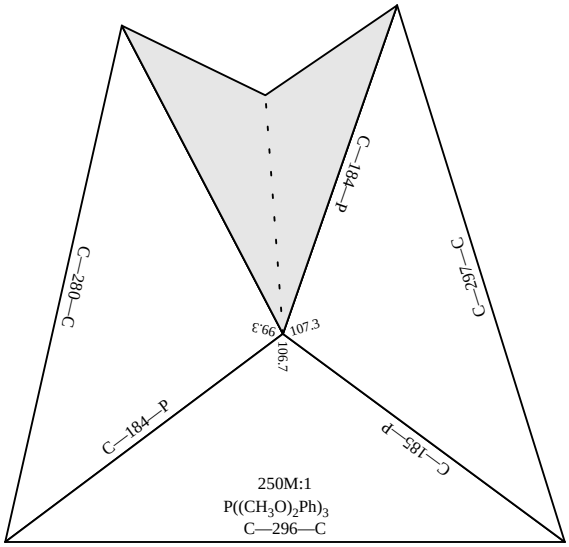
Molecular Origami of P((CH3O)2Ph)3

!P1
C1
C17
C9
P((CH3O)2Ph)3

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: 187 310 417 531 actual: 197 320 407 521 center: 302 420 actual size: 210 201
Better: (centerx 5.71) (centery 5.56) (scale 250)
%%BoundingBox: 288 350 518 571 actual: 298 360 508 561 center: 403 461 actual size: 210 201