

# Molecular Origami of BiPh3

given information

ElementNames	[ (Bi) (C) (C) (C) ]	
dotted	F	
distance	221.114	Bi <sup>1</sup> -C <sup>1</sup>
distance	225.051	Bi <sup>1</sup> -C <sup>16</sup>
distance	225.191	Bi <sup>1</sup> -C <sup>5</sup>
angle	91.969	C <sup>16</sup> -Bi <sup>1</sup> -C <sup>1</sup>
	320.9	C <sup>16</sup> -C <sup>1</sup>
angle	94.228	C <sup>5</sup> -Bi <sup>1</sup> -C <sup>16</sup>
	329.9	C <sup>5</sup> -C <sup>16</sup>
angle	95.614	C <sup>5</sup> -Bi <sup>1</sup> -C <sup>1</sup>
	330.7	C <sup>5</sup> -C <sup>1</sup>
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

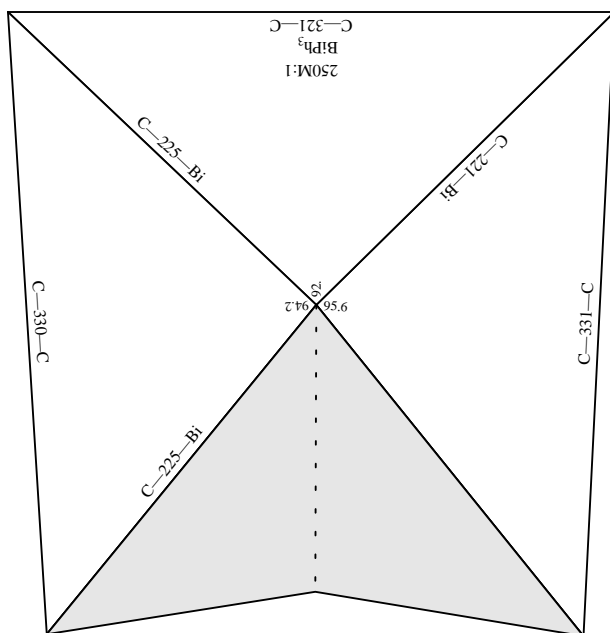
# Molecular Origami of BiPh3

!Bi1  
C1  
C16  
C5  
BiPh3

trigonal pyramidal

scale 250,000,000 : 1  
units: pm  
offsetx -1.67 offsety 0.9

View -1



Current: (centerx 2.63) (centery 5.90) (scale 250)

%%BoundingBox: 179 294 427 548 actual: 189 304 417 538

center: 303 421

actual size: 227 234

Better: (centerx 2.67) (centery 5.55) (scale 250)

%%BoundingBox: 62 334 309 588

actual: 72 344 299 578

center: 186 461

actual size: 227 234