Molecular Origami of Bi((CH3)3Ph)3

given information

ElementNames	[(Bi) (C) (C) (C)]	
dotted	F	
distance	224.332	Bi^{1}-C^{10}
distance	231.285	Bi^{1}-C^{19}
distance	231.550	Bi^{1}-C^{1}
angle	95.201	C^{19}-Bi^{1}-C^{1}
	341.8	C^{19}-C^{1}
angle	105.526	C^{19}-Bi^{1}-C^{10}
	362.8	C^{19}-C^{10}
angle	107.413	C^{10}-Bi^{1}-C^{1}
	367.5	C^{10}-C^{1}
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

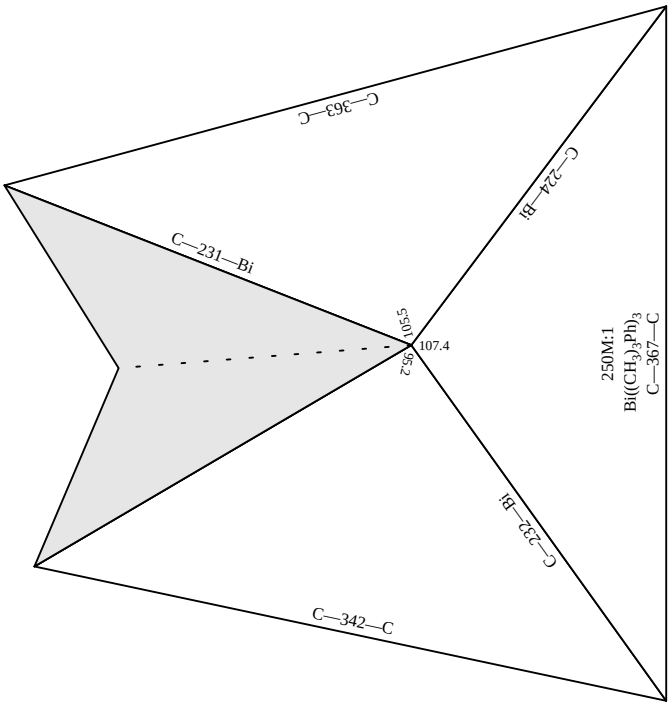
Molecular Origami of Bi((CH3)3Ph)3

!Bi1
C1
C10
C19
Bi((CH3)3Ph)3

trigonal pyramidal

scale 250,000,000 : 1
units: pm
offsetx -0.05 offsety 2.67

View -1



Current: (centerx 4.25) (centery 7.67) (scale 250)
%%BoundingBox: 178 282 446 562 actual: 188 292 436 552 center: 312 422 actual size: 248 260
Better: (centerx 4.16) (centery 7.31) (scale 250)
%%BoundingBox: 168 448 436 728 actual: 178 458 426 718 center: 302 588 actual size: 248 260