

# Molecular Origami of Bi((EtO)2Ph)3

given information

ElementNames	[ (Bi) (C) (C) (C) ]	
dotted	F	
distance	225.414	Bi <sup>1</sup> -C <sup>1</sup>
distance	225.621	Bi <sup>1</sup> -C <sup>13</sup>
distance	227.784	Bi <sup>1</sup> -C <sup>7</sup>
angle	95.683	C <sup>13</sup> -Bi <sup>1</sup> -C <sup>1</sup>
	334.3	C <sup>13</sup> -C <sup>1</sup>
angle	101.174	C <sup>7</sup> -Bi <sup>1</sup> -C <sup>13</sup>
	350.3	C <sup>7</sup> -C <sup>13</sup>
angle	101.373	C <sup>7</sup> -Bi <sup>1</sup> -C <sup>1</sup>
	350.6	C <sup>7</sup> -C <sup>1</sup>
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

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!Bi1

C1

C13

C7

Bi((EtO)2Ph)3

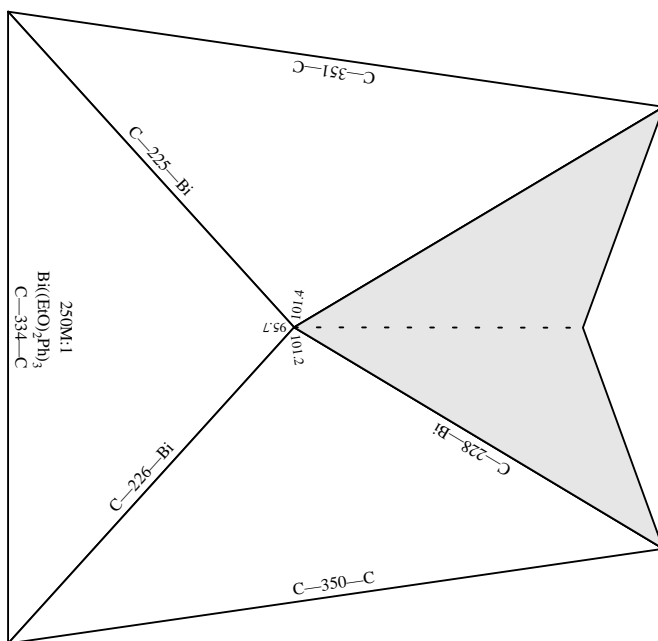
trigonal pyramidal

scale 250,000,000 : 1

units: pm

offsetx -0.05 offsety -0.81

View -1



Current: (centerx 4.25) (centery 4.19) (scale 250)

%%BoundingBox: 178 292 443 549 actual: 188 302 433 539

center: 310 420

actual size: 246 237

Better: (centerx 4.19) (centery 3.85) (scale 250)

%%BoundingBox: 169 209 435 466

actual: 179 219 425 456

center: 302 338

actual size: 246 237