

# Molecular Origami of N(CH3)3

given information

ElementNames	[ (N) (C) (C) (C) ]	
dotted	F	
distance	144.778	N <sup>1</sup> -C <sup>1</sup>
distance	144.804	N <sup>1</sup> -C <sup>2</sup>
distance	144.835	N <sup>1</sup> -C <sup>3</sup>
angle	110.628	C <sup>3</sup> -N <sup>1</sup> -C <sup>2</sup>
	238.2	C <sup>3</sup> -C <sup>2</sup>
angle	110.647	C <sup>3</sup> -N <sup>1</sup> -C <sup>1</sup>
	238.2	C <sup>3</sup> -C <sup>1</sup>
angle	110.657	C <sup>2</sup> -N <sup>1</sup> -C <sup>1</sup>
	238.2	C <sup>2</sup> -C <sup>1</sup>
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XABC

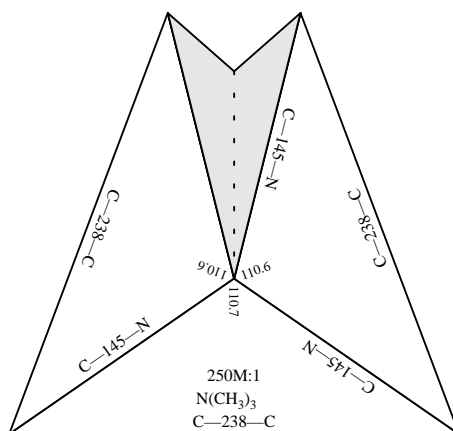
# Molecular Origami of N(CH3)3

!N1  
C1  
C2  
C3  
N(CH3)3

trigonal pyramidal

scale 250,000,000 : 1  
units: pm  
offsetx 1.05 offsety 0.9

View -1



Current: (centerx 5.35) (centery 5.90) (scale 250)

%%BoundingBox: 206 330 395 508 actual: 216 340 385 498

Better: (centerx 5.42) (centery 5.58) (scale 250)

%%BoundingBox: 287 372 476 550 actual: 297 382 466 540

center: 301 419

actual size: 169 158

center: 382 461

actual size: 169 158