Molecular Origami of NEt3(CH3)^+

given information

ElementNames	[(N) (C) (C) (C) (C)]	
distance	150.227	N^1-C^4
distance	150.980	N^1-C^1
distance	151.511	N^1-C^2
distance	151.694	N^1-C^3
angle	106.598	C^4-N^1-C^1
	241.5	C^4-C^1
angle	108.788	C^3-N^1-C^2
	246.5	C^3-C^2
angle	109.247	C^3-N^1-C^1
	246.8	C^3-C^1
angle	109.723	C^4-N^1-C^2
	246.8	C^4-C^2
angle	110.185	C^4-N^1-C^3
	247.6	C^4-C^3
angle	112.282	C^2-N^1-C^1
	251.2	C^2-C^1
dopage	T	
AutoAlign	F	

structure type: XABCD

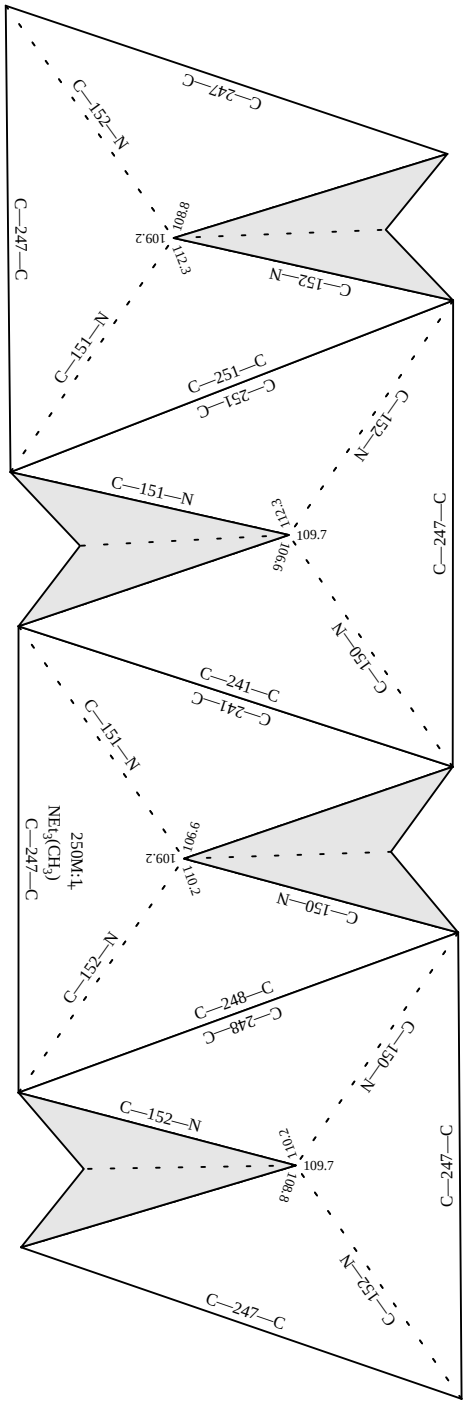
Molecular Origami of NEt3(CH3)^+

!N1
C1
C2
C3
C4
NEt3(CH3)^+

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.31 offsety 0.08

View -1



Current: (centerx 3.99) (centery 5.08) (scale 250)
%%BoundingBox: 210 153 401 696 actual: 220 163 391 686 center: 306 424 actual size: 171 522
Better: (centerx 3.99) (centery 4.69) (scale 250)
%%BoundingBox: 188 131 379 673 actual: 198 141 369 663 center: 284 402 actual size: 171 522