Molecular Origami of CH3CN

given information

ElementNames	[(C) (C) (N)]	
dotted	F	
distance	113.190	C^1-N^1
distance	145.429	C^1-C^2
angle	178.897	N^1-C^1-C^2
	258.6	N^1-C^2
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

Molecular Origami of CH3CN

!C1
C2
N1
CH3CN

linear

scale 250,000,000 : 1
units: pm
offsetx -0.08 offsety 0.9

View -1

Current: (centerx 4.22) (centery 5.90) (scale 250)
%%BoundingBox: 294 128 314 435 actual: 304 138 304 425 center: 304 282 actual size: 0 286
Better: (centerx 4.25) (centery 7.49) (scale 250)
%%BoundingBox: 290 308 310 614 actual: 300 318 300 604 center: 300 461 actual size: 0 286