

# Molecular Origami of CH3CN

given information

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

C<sup>1</sup>-N<sup>1</sup>  
C<sup>1</sup>-C<sup>2</sup>  
N<sup>1</sup>-C<sup>1</sup>-C<sup>2</sup>  
N<sup>1</sup>-C<sup>2</sup>

structure type: XAB

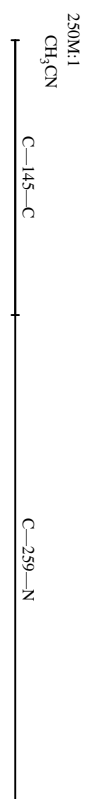
# Molecular Origami of CH<sub>3</sub>CN

!C1  
C2  
N1  
CH<sub>3</sub>CN

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