Molecular Origami of CN^-

given information

ElementNames	[(C) (N)]	
distance	92.258	C^{1}-N^{1}
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XA

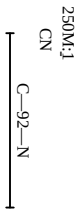
Molecular Origami of CN^-

!C1
N1
CN^-

diatomic

scale 250,000,000 : 1
units: pm
offsetx -0.05 offsety 0.9

View -1



Current: (centerx 4.25) (centery 5.90) (scale 250)
%%BoundingBox: 296 349 316 435 actual: 306 359 306 425 center: 306 392 actual size: 0 65
Better: (centerx 4.25) (centery 5.95) (scale 250)
%%BoundingBox: 292 418 312 503 actual: 302 428 302 493 center: 302 461 actual size: 0 65