Molecular Origami of NCO^-

given information

ElementNames	[(C) (N) (O)]	
dotted	F	
distance	114.794	C^1-N^1
distance	121.574	C^1-O^1
angle	179.806	O^1-C^1-N^1
	236.4	O^1-N^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

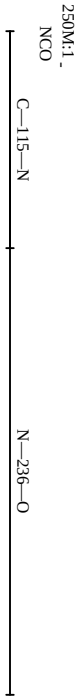
Molecular Origami of NCO^-

!C1
N1
O1
NCO^-

linear

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: 296 166 316 435 actual: 306 176 306 425 center: 306 300 actual size: 0 249
Better: (centerx 4.25) (centery 7.23) (scale 250)
%%BoundingBox: 292 326 312 595 actual: 302 336 302 585 center: 302 461 actual size: 0 249