

Molecular Origami of NCO^-
given information

ElementNames	[(C) (N) (O)]
dotted	F
distance	112.730
distance	121.992
angle	179.062
	234.7
dopage	T
AutoAlign	F
showboth	F

C¹-O¹
C¹-N¹
O¹-C¹-N¹
O¹-N¹

structure type: XAB

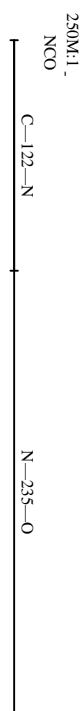
Molecular Origami of NCO⁻

!C1
N1
O1
NCO⁻

linear

scale 250,000,000 : 1
units: pm
offsetx -0.07 offsety 0.9

View -1



Current: (centerx 4.23) (centery 5.90) (scale 250)

%%BoundingBox: 295 162 315 435 actual: 305 172 305 425

Better: (centerx 4.25) (centery 7.26) (scale 250)

%%BoundingBox: 291 324 311 597 actual: 301 334 301 587

center: 305 298

actual size: 0 253

center: 301 461

actual size: 0 253