Molecular Origami of HNCO

given information

ElementNames	[(C) (N) (O)]	
dotted	F	
distance	120.229	C^1-O^1
distance	123.654	C^1-N^1
angle	178.155	O^1-C^1-N^1
	243.9	O^1-N^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

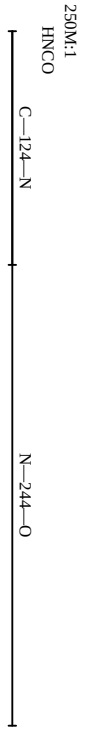
Molecular Origami of HNCO

!C1
N1
O1
HNCO

linear

scale 250,000,000 : 1
units: pm
offsetx -0.09 offsety 0.9

View -1



Current: (centerx 4.21) (centery 5.90) (scale 250)
%%BoundingBox: 293 154 313 435 actual: 303 164 303 425 center: 303 295 actual size: 0 260
Better: (centerx 4.25) (centery 7.31) (scale 250)
%%BoundingBox: 290 321 310 601 actual: 300 331 300 591 center: 300 461 actual size: 0 260