Molecular Origami of N2F2trans

given information

ElementNames	[(N) (F) (N)]	
dotted	F	
distance	123.004	N^1-N^2
distance	139.995	N^1-F^1
angle	105.995	N^2-N^1-F^1
	210.3	N^2-F^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

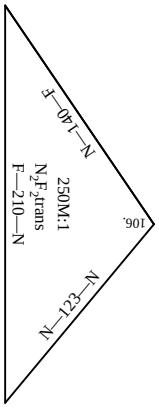
Molecular Origami of N2F2trans

!N1
F1
N2
N2F2trans

bent

scale 250,000,000 : 1
units: pm
offsetx -1.07 offsety 0.9

View -1



Current: (centerx 3.23) (centery 5.90) (scale 250)
%%BoundingBox: 223 340 298 509 actual: 233 350 288 499 center: 260 425 actual size: 56 149
Better: (centerx 3.86) (centery 5.50) (scale 250)
%%BoundingBox: 191 376 267 545 actual: 201 386 257 535 center: 229 461 actual size: 56 149