Molecular Origami of N2F2cis

given information

ElementNames	[(N) (F) (N)]	
dotted	F	
distance	120.997	N^1-N^2
distance	138.006	N^1-F^1
angle	114.502	N^2-N^1-F^1
	218.	N^2-F^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

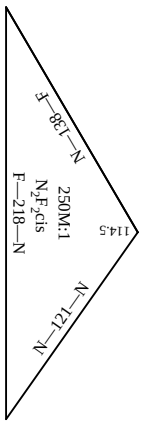
Molecular Origami of N2F2cis

!N1
F1
N2
N2F2cis

bent

scale 250,000,000 : 1
units: pm
offsetx -1.06 offsety 0.9

View -1



Current: (centerx 3.24) (centery 5.90) (scale 250)
%%BoundingBox: 223 338 293 512 actual: 233 348 283 502 center: 258 425 actual size: 49 155
Better: (centerx 3.91) (centery 5.50) (scale 250)
%%BoundingBox: 195 374 264 548 actual: 205 384 254 538 center: 230 461 actual size: 49 155