

Molecular Origami of N2H2

given information

ElementNames	[(N) (H) (N)]
dotted	F
distance	101.403
distance	123.003
angle	99.999
	172.5
dopage	T
AutoAlign	F
showboth	F

N¹-H¹
N¹-N²
N²-N¹-H¹
N²-H¹

structure type: XAB

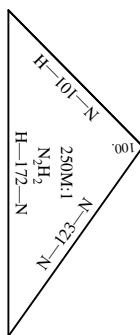
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!N1
H1
N2
N2H2

bent

scale 250,000,000 : 1
units: pm
offsetx -0.85 offsety 0.9

View -1



Current: (centerx 3.45) (centery 5.90) (scale 250)

%%BoundingBox: 238 354 309 496 actual: 248 364 299 486

center: 274 425

actual size: 50 122

Better: (centerx 3.90) (centery 5.50) (scale 250)

%%BoundingBox: 210 390 280 532

actual: 220 400 270 522

center: 245 461

actual size: 50 122