Molecular Origami of I3^-

given information

ElementNames	[(I) (I) (I)]	
dotted	F	
distance	283.008	I^2-I^3
distance	303.830	I^2-I^1
angle	173.701	I^3-I^2-I^1
	586.	I^3-I^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

Molecular Origami of I3^-

!I2
I1
I3
I3^-

linear

scale 250,000,000 : 1
units: pm
offsetx -0.38 offsety 0.9

View -1

Current: (centerx 3.92) (centery 5.90) (scale 250)
%%BoundingBox: 272 -216 292 435 actual: 282 -206 282 425 center: 282 110 actual size: 0 631
Better: (centerx 4.25) (centery 9.88) (scale 250)
%%BoundingBox: 269 136 289 786 actual: 279 146 279 776 center: 279 461 actual size: 0 631