Molecular Origami of I3^-

given information

ElementNames	[(I) (I) (I)]	
dotted	F	
distance	292.586	I^1-I^3
distance	292.586	I^1-I^2
angle	180.000	I^3-I^1-I^2
	585.2	I^3-I^2
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

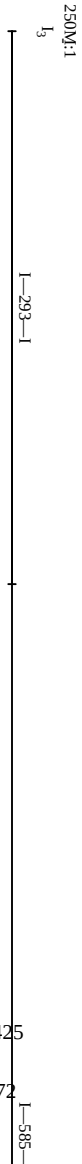
Molecular Origami of I3^-

!I1
I2
I3
I3^-

linear

scale 250,000,000 : 1
units: pm
offsetx -0.05 offsety 0.9

View -1



Current: (centerx 4.25) (centery 5.90) (scale 250)
%%BoundingBox: 296 -207 316 435 actual: 306 -197 306 425 center: 306 114 actual size: 0 622
Better: (centerx 4.25) (centery 9.82) (scale 250)
%%BoundingBox: 292 140 312 782 actual: 302 150 302 772 center: 302 461 actual size: 0 622