Molecular Origami of BrIBr^-

given information

ElementNames	[(I) (Br) (Br)]	
dotted	F	
distance	262.602	I^1-Br^1
distance	278.275	I^1-Br^2
angle	178.207	Br^2-I^1-Br^1
	540.8	Br^2-Br^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

Molecular Origami of BrIBr^-

!I1
Br1
Br2
BrIBr^-

linear

scale 250,000,000 : 1
units: pm
offsetx -0.13 offsety 0.9

View -1

Current: (centerx 4.17) (centery 5.90) (scale 250)
%%BoundingBox: 290 -155 310 435 actual: 300 -145 300 425 center: 300 140 actual size: 0 569
Better: (centerx 4.25) (centery 9.45) (scale 250)
%%BoundingBox: 287 166 307 755 actual: 297 176 297 745 center: 297 461 actual size: 0 569