Molecular Origami of BrIBr^-

given information

ElementNames	[(I) (Br) (Br)]	
dotted	F	
distance	268.518	I^1-Br^1
distance	271.807	I^1-Br^2
angle	179.740	Br^2-I^1-Br^1
	540.3	Br^2-Br^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

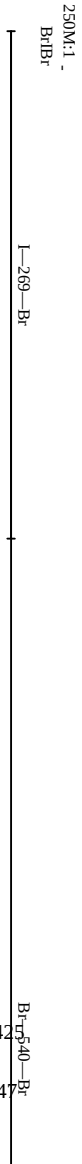
Molecular Origami of BrIBr^-

!I1
Br1
Br2
BrIBr^-

linear

scale 250,000,000 : 1
units: pm
offsetx -0.06 offsety 0.9

View -1



Current: (centerx 4.24) (centery 5.90) (scale 250)
%%BoundingBox: 295 -158 315 435 actual: 305 -148 305 423 center: 305 138 actual size: 0 573
Better: (centerx 4.25) (centery 9.48) (scale 250)
%%BoundingBox: 292 164 312 757 actual: 302 174 302 747 center: 302 461 actual size: 0 573