Molecular Origami of Br3^-

given information

ElementNames	[(Br) (Br) (Br)]	
dotted	F	
distance	255.100	Br^1-Br^2
distance	255.110	Br^1-Br^3
angle	177.610	Br^3-Br^1-Br^2
	510.1	Br^3-Br^2
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

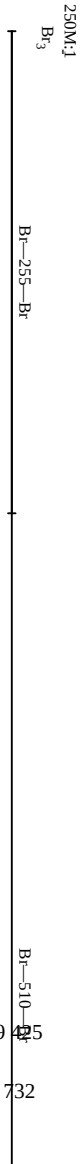
Molecular Origami of Br3^-

!Br1
Br2
Br3
Br3^-

linear

scale 250,000,000 : 1
units: pm
offsetx -0.15 offsety 0.9

View -1



Current: (centerx 4.15) (centery 5.90) (scale 250)
%%BoundingBox: 289 -127 309 435 actual: 299 -117 299 425 center: 299 154 actual size: 0 542
Better: (centerx 4.25) (centery 9.27) (scale 250)
%%BoundingBox: 285 180 305 742 actual: 295 190 295 732 center: 295 461 actual size: 0 542