Molecular Origami of ClICl^-

given information

ElementNames	[(I) (Cl) (Cl)]	
dotted	F	
distance	253.982	I^1-Cl^2
distance	253.982	I^1-Cl^1
angle	179.999	Cl^2-I^1-Cl^1
	508.	Cl^2-Cl^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

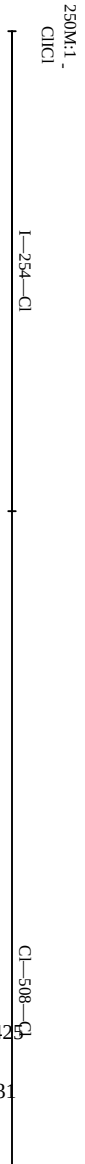
Molecular Origami of ClICl^-

!I1
Cl1
Cl2
ClICl^-

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 -125 316 435 actual: 306 -115 306 425 center: 306 155 actual size: 0 540
Better: (centerx 4.25) (centery 9.25) (scale 250)
%%BoundingBox: 292 181 312 741 actual: 302 191 302 731 center: 302 461 actual size: 0 540