Molecular Origami of ClICl^-

given information

ElementNames	[(I) (Cl) (Cl)]	
dotted	F	
distance	257.413	I^1-Cl^2
distance	257.413	I^1-Cl^1
angle	179.999	Cl^2-I^1-Cl^1
	514.8	Cl^2-Cl^1
dopage	T	
AutoAlign	F	
showboth	F	

structure type: XAB

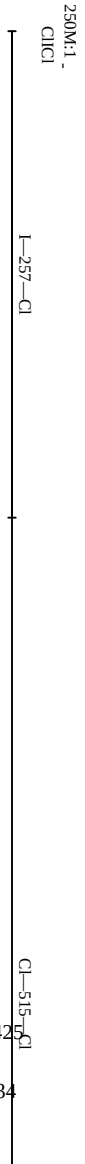
Molecular Origami of ClICl^-

!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 -132 316 435 actual: 306 -122 306 425 center: 306 151 actual size: 0 547
Better: (centerx 4.25) (centery 9.30) (scale 250)
%%BoundingBox: 292 177 312 744 actual: 302 187 302 734 center: 302 461 actual size: 0 547