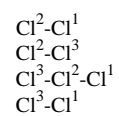


# Molecular Origami of Cl3^-

given information

ElementNames	[ (Cl) (Cl) (Cl) ]
dotted	F
distance	224.761
distance	233.811
angle	177.543
	458.5
dopage	T
AutoAlign	F
showboth	F



structure type: XAB

Molecular Origami of Cl3^-

!Cl2  
Cl1  
Cl3  
Cl3^-

linear

scale 250,000,000 : 1  
units: pm  
offsetx -0.14 offsety 0.9

View -1



Current: (centerx 4.16) (centery 5.90) (scale 250)			
%%BoundingBox: 290 -69 310 435	actual: 300 -59 300 425	center: 300 183	actual size: 0 484
Better: (centerx 4.25) (centery 8.86) (scale 250)			
%%BoundingBox: 286 209 306 713	actual: 296 219 296 703	center: 296 461	actual size: 0 484