Molecular Origami of CI4

given information

ElementNames	[(C) (I) (I) (I) (I)]	
distance	215.439	C^1-I^2
distance	215.439	C^1-I^3
distance	215.439	C^1-I^4
distance	215.439	C^1-I^1
angle	109.245	I^4-C^1-I^2
	351.3	I^4-I^2
angle	109.245	I^3-C^1-I^1
	351.3	I^3-I^1
angle	109.584	I^4-C^1-I^3
	352.1	I^4-I^3
angle	109.584	I^3-C^1-I^2
	352.1	I^3-I^2
angle	109.584	I^4-C^1-I^1
	352.1	I^4-I^1
angle	109.584	I^2-C^1-I^1
	352.1	I^2-I^1
dopage	T	
AutoAlign	F	

structure type: XABCD

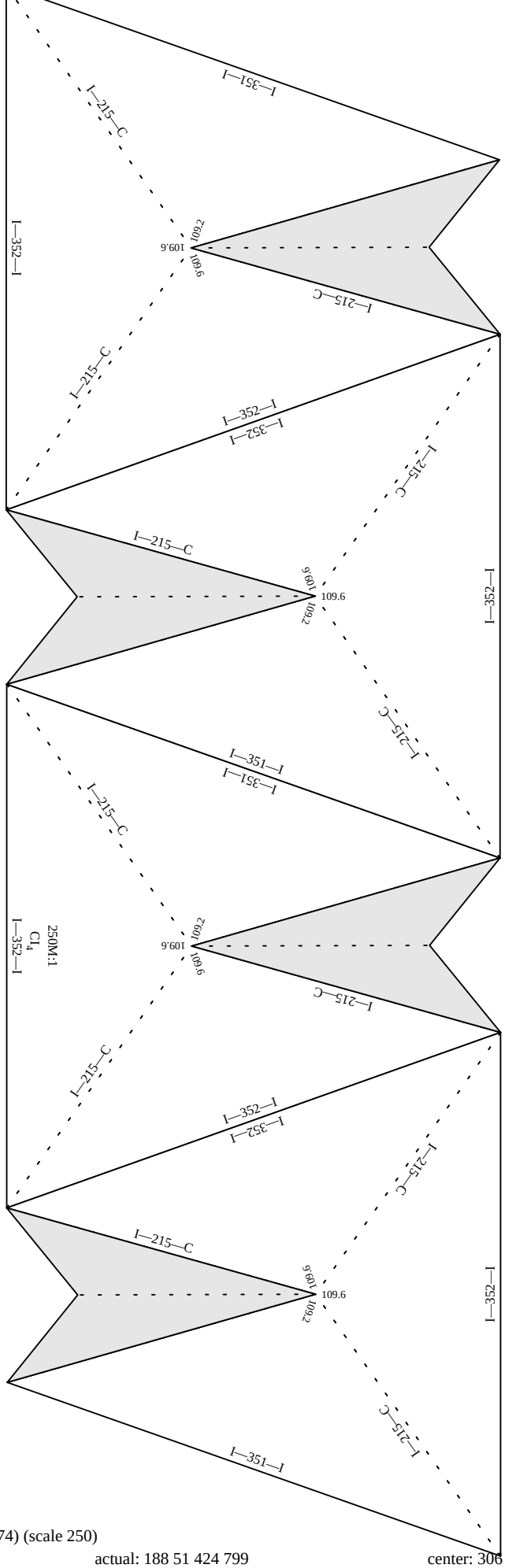
Molecular Origami of CI4

!C1
I1
I2
I3
I4
CI4

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.46 offsety -0.26

View -1



Current: (centerx 3.84) (centery 4.74) (scale 250)
%%BoundingBox: 178 41 434 809 actual: 188 51 424 799 center: 306 425 actual size: 235 748
Better: (centerx 4.00) (centery 4.48) (scale 225)
%%BoundingBox: 157 17 389 738 actual: 167 27 379 728 center: 273 377 actual size: 212 701