Molecular Origami of CH2I2

given information

ElementNames	[(C) (H) (H) (I) (I)]	
distance	110.427	C^1-H^2
distance	110.427	C^1-H^1
distance	213.660	C^1-I^2
distance	213.660	C^1-I^1
angle	108.372	I^2-C^1-H^1
	269.7	I^2-H^1
angle	108.372	I^2-C^1-H^2
	269.7	I^2-H^2
angle	108.372	I^1-C^1-H^2
	269.7	I^1-H^2
angle	108.372	I^1-C^1-H^1
	269.7	I^1-H^1
angle	110.441	H^2-C^1-H^1
	181.4	H^2-H^1
angle	112.914	I^2-C^1-I^1
	356.2	I^2-I^1
dopage	T	
AutoAlign	F	

structure type: XABCD

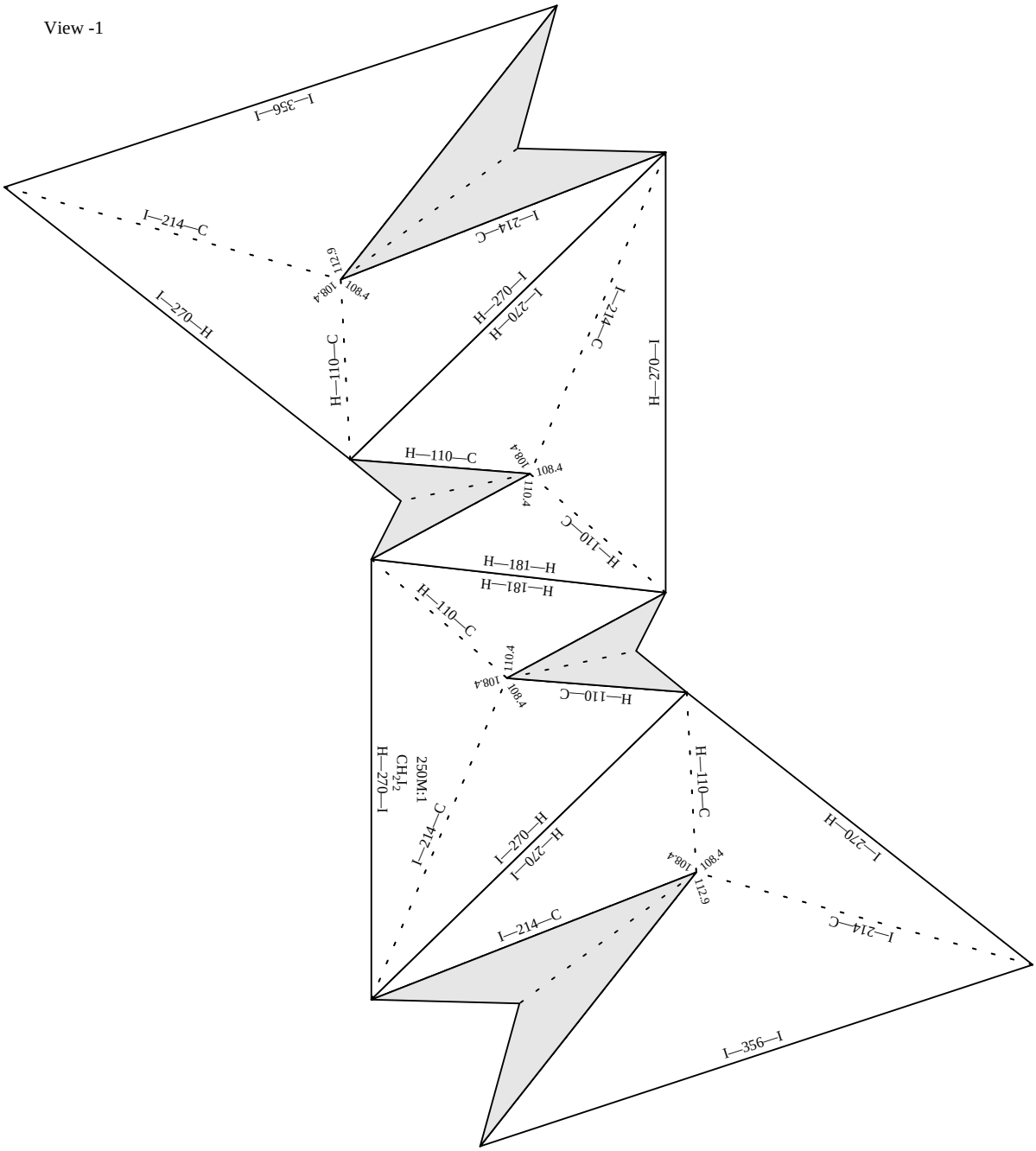
Molecular Origami of CH2I2

!C1
H1
H2
I1
I2
CH2I2

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.51 offsety -0.65

View -1



Current: (centerx 3.79) (centery 4.35) (scale 250)
%%BoundingBox: 45 100 511 615 actual: 55 110 501 605 center: 278 358 actual size: 446 495
Better: (centerx 4.18) (centery 4.88) (scale 250)
%%BoundingBox: 36 92 502 607 actual: 46 102 492 597 center: 269 349 actual size: 446 495