Molecular Origami of CH2I2

given information

ElementNames	[(C) (H) (H) (I) (I)]	
distance	99.997	C^1-H^1
distance	100.003	C^1-H^2
distance	212.597	C^1-I^1
distance	212.605	C^1-I^2
angle	108.247	I^2-C^1-H^1
	261.8	I^2-H^1
angle	108.249	I^1-C^1-H^2
	261.8	I^1-H^2
angle	108.252	I^1-C^1-H^1
	261.8	I^1-H^1
angle	108.253	I^2-C^1-H^2
	261.8	I^2-H^2
angle	109.500	H^2-C^1-H^1
	163.3	H^2-H^1
angle	114.277	I^2-C^1-I^1
	357.2	I^2-I^1
dopage	T	
AutoAlign	F	

structure type: XABCD

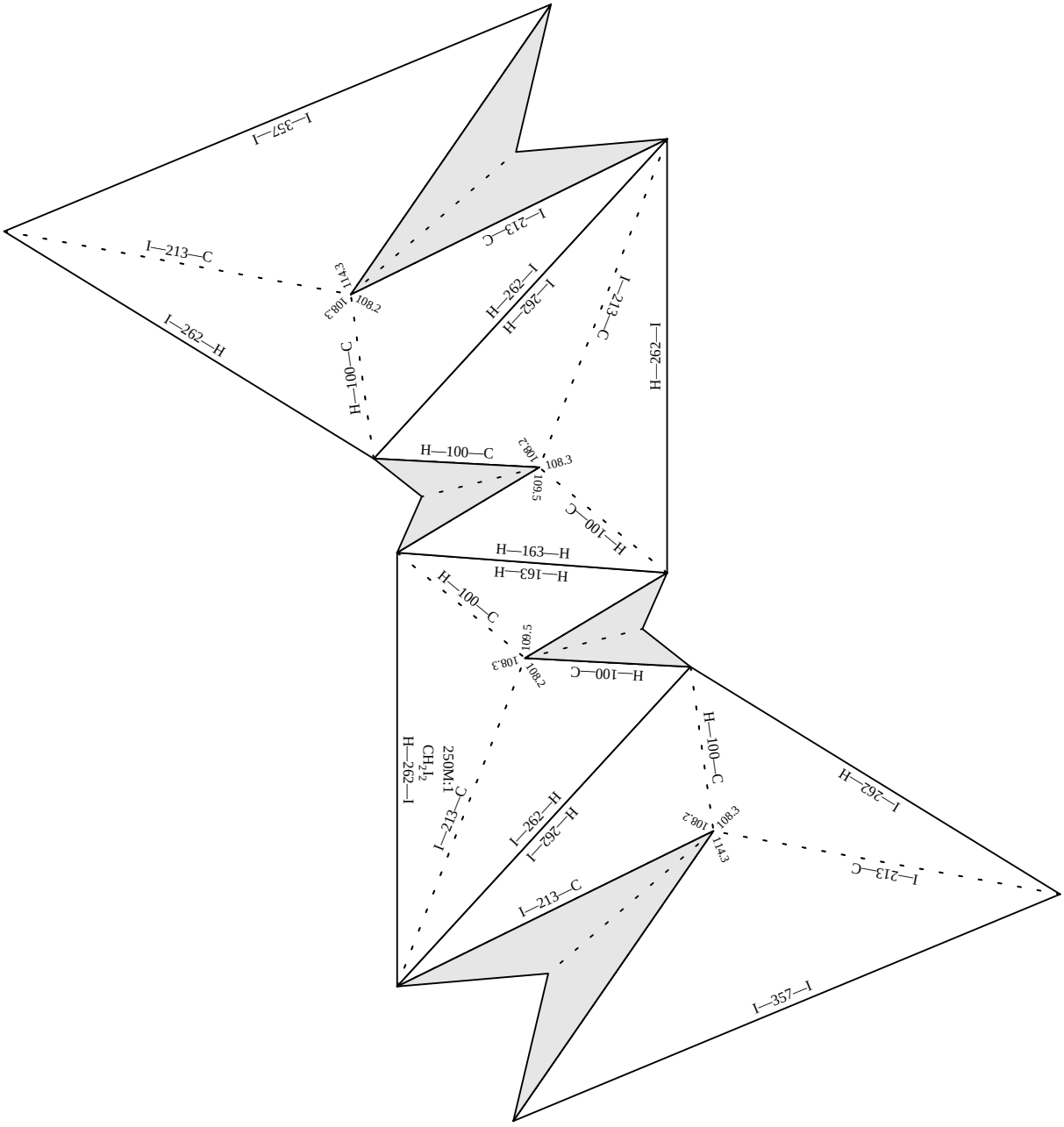
Molecular Origami of CH2I2

!C1
H1
H2
I1
I2
CH2I2

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.49 offsety -0.7

View -1



Current: (centerx 3.81) (centery 4.30) (scale 250)
%%BoundingBox: 42 102 513 599 actual: 52 112 503 589 center: 277 350 actual size: 451 477
Better: (centerx 4.21) (centery 4.93) (scale 250)
%%BoundingBox: 35 97 506 594 actual: 45 107 496 584 center: 271 346 actual size: 451 477