Molecular Origami of CH2F2

given information

ElementNames	[(C) (F) (F) (H) (H)]	
distance	107.941	C^1-H^2
distance	108.030	C^1-H^1
distance	128.676	C^1-F^1
distance	141.083	C^1-F^2
angle	99.893	F^2-C^1-F^1
	206.6	F^2-F^1
angle	105.022	H^2-C^1-F^1
	188.2	H^2-F^1
angle	109.495	H^2-C^1-H^1
	176.4	H^2-H^1
angle	113.070	H^1-C^1-F^2
	208.6	H^1-F^2
angle	113.659	H^2-C^1-F^2
	209.2	H^2-F^2
angle	115.268	H^1-C^1-F^1
	200.2	H^1-F^1
dopage	T	
AutoAlign	F	

structure type: XABCD

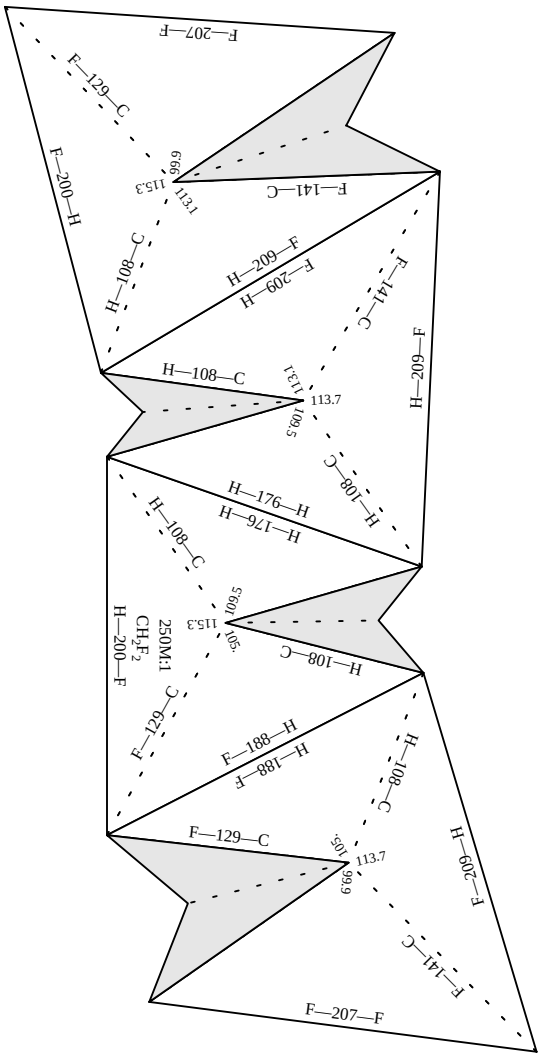
Molecular Origami of CH2F2

!C1
F1
F2
H1
H2
CH2F2

special tetrahedral

scale 250,000,000 : 1
units: pm
offsetx -0.38 offsety 0.24

View -1



Current: (centerx 3.92) (centery 5.24) (scale 250)
%%BoundingBox: 189 206 409 618 actual: 199 216 399 608 center: 299 412 actual size: 199 392
Better: (centerx 4.02) (centery 5.01) (scale 250)
%%BoundingBox: 169 207 388 619 actual: 179 217 378 609 center: 279 413 actual size: 199 392