

Molecular Origami of CHI3

given information

ElementNames	[(C) (H) (I) (I) (I)]	
distance	99.999	C ¹ -H ¹
distance	210.322	C ¹ -I ³
distance	214.724	C ¹ -I ¹
distance	221.964	C ¹ -I ²
angle	104.491	I ¹ -C ¹ -H ¹
	258.6	I ¹ -H ¹
angle	107.906	I ³ -C ¹ -I ²
	349.6	I ³ -I ²
angle	108.927	I ³ -C ¹ -H ¹
	260.5	I ³ -H ¹
angle	110.168	I ² -C ¹ -H ¹
	273.1	I ² -H ¹
angle	112.060	I ² -C ¹ -I ¹
	362.2	I ² -I ¹
angle	113.226	I ³ -C ¹ -I ¹
	354.9	I ³ -I ¹
dopage	T	
AutoAlign	F	

structure type: XABCD

!C1
H1
I1
I2
I3
CHI3

scale 250,000,000 : 1

```
offsetx -0.01  offsety -0.35
```

[illegible]

```
%%BoundingBox: 29 79 527 625
```

actual: 39 89 517 615

center: 278 352

actual size: 478 525

Better: (centerx 4.68) (centery 5.26) (scale 250)

%%BoundingBox: 57 98 554 643

actual: 67 108 544 633

center: 305 371

actual size: 478 525