

Molecular Origami of CHI3

given information

ElementNames	[(C) (H) (I) (I) (I)]	
distance	100.000	C ¹ -H ¹
distance	214.084	C ¹ -I ³
distance	214.084	C ¹ -I ¹
distance	220.413	C ¹ -I ²
angle	108.101	I ³ -C ¹ -H ¹
	262.9	I ³ -H ¹
angle	108.101	I ¹ -C ¹ -H ¹
	262.9	I ¹ -H ¹
angle	109.538	I ³ -C ¹ -I ²
	354.9	I ³ -I ²
angle	109.538	I ² -C ¹ -I ¹
	354.9	I ² -I ¹
angle	110.075	I ² -C ¹ -H ¹
	271.5	I ² -H ¹
angle	111.462	I ³ -C ¹ -I ¹
	353.8	I ³ -I ¹
dopage	T	
AutoAlign	F	

structure type: XABCD

!C1
H1
I1
I2
I3
CHI3

```
scale 250,000,000 : 1
units: pm
offsetx -0.07 offsety -0.46
```

[illegible]

actual size: 492 523

actual size: 492 523