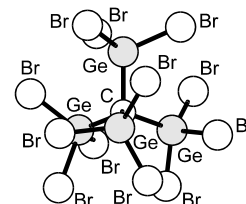


16 **CBr₁₂Ge₄**ED, DFT
calculations**Methanetetrayltetrakis[tribromogermene]****T assumed**
C(GeBr₃)₄

r_a	$\text{\AA}^a)$	θ_a	$\text{deg}^a)$
C–Ge	2.042(8)	C–Ge–Br	112.9(5)
Ge–Br	2.282(3)	Br–Ge–Br	105.9(5)
		$\tau^b)$	31.4(9)

The twelve Br atom positions are close to the vertices of a regular icosahedron. BPW91/LanL2DZ calculations closely reproduced the observed structure.

The nozzle temperature was 218 °C.



^{a)} Twice the estimated standard errors including a systematic error.

^{b)} Ge–C–Ge–Br torsional angle from the *syn* position.

Haaland, A., Shorokhov, D.J., Strand, T.G., Kouvetakis, J., O’Keeffe, M.: Inorg. Chem. **36** (1997) 5198.