

Structure Data of Free Polyatomic Molecules

1 MW	C Ag BrO		Bromocarbonylsilver	C_{Oov} Br–Ag–C=O
	r_0	\AA^{a}		
	C=O	1.1236(9)		
	C–Ag	2.027(1)		
	Ag–Br	2.3731(5)		
	r_{IE}	\AA^{a}		
	C=O	1.1242(1)		
	C–Ag	2.0275(1)		
	Ag–Br	2.3704(2)		
	$r_{\text{m}}^{(1)}$	\AA^{a}		
	C=O	1.1236(1)		
	C–Ag	2.0264(2)		
	Ag–Br	2.3692(2)		
	$r_{\text{m}}^{(2)}$	\AA^{a}		
	C=O	1.1219(4)		
	C–Ag	2.02800(4)		
	Ag–Br	2.36826(2)		

^a) Estimated standard errors.

Walker, N.R., Gerry, M.C.L.: Inorg. Chem. **41** (2002) 1236.