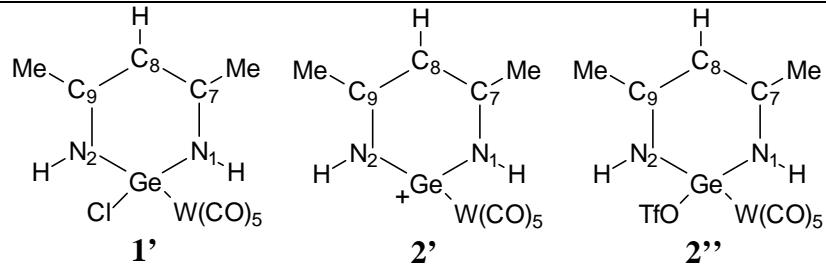


¹H-, ¹³C-{¹H}- and ¹⁹F-{¹H} NMR (C₆D₆) and IR data for compounds 2, 3, 13

	2	3	13
¹⁹ F NMR (δ , ppm)			
CF ₃ SO ₃	- 1.65	- 2.03	-
¹ H NMR (δ , ppm)			
CH	5.24, 5.85 ^a	4.98, 5.66 ^a	4.96, 5.38 ^a
CH ₃	1.58, 2.11 ^a	1.44, 2.07 ^a	1.45, 1.88 ^a
¹³ C NMR (δ , ppm)			
CH	103.8	103.5	101.1 ^a
CH ₃	24.0	23.9	24.8 ^a
CO	194.6, 197.4	-	201.0 ^a
IR : v _{CO} (cm ⁻¹);			
v _{OTf} (cm ⁻¹)			
CHCl ₃	2062.4, 1973.9, 1932.3; 1973.1366.5	-	1897.8
Pyridine	1366.5	1378.7	
C ₆ D ₆	2063.8, 1934.4, 1920.5; 1379.1, 1274.0 2079.3, 1986.2, 1943.1; 1365.4	1367.4, 1271.7	

^a: CDCl₃

Calculated Geometrical Parameters of **1', **2'** and **2''***** (Bond Length in Angstroms and Bond angles in degrees).



GeCl	2.359	-	-
GeO	-	-	2.027
GeN ₁	1.950	1.902	1.927
GeN ₂	1.950	1.902	1.935
N ₁ C ₇	1.353	1.367	1.358
C ₇ C ₈	1.411	1.408	1.406
C ₉ C ₈	1.410	1.408	1.413
N ₂ C ₉	1.353	1.367	1.351
GeW	2.637	2.538	2.630
N ₁ GeN ₂	86.84	91.8	91.6
N ₁ GeCl	97.7	-	-
N ₂ GeCl	97.7	-	-
N ₁ GeO	-	-	90.7
N ₂ GeO	-	-	94.1
WGeCl	121.7	-	-
WGeO	-	-	120.0
N ₁ GeW	121.1	134.1	123.1
N ₂ GeW	121.2	134.1	127.8

* In the case of **2''**, the frequencies dealing with the total movement of the molecule has a very low value.

Total Natural Charge of 1', 2' and 2'' (NBO calculations).

	1'	2'	2''
Ge	1.375	1.486	1.574
N ₁	-0.951	-0.978	-0.968
N ₂	-0.951	-0.979	-0.956
Cl	-0.499	-	-
O	-	-	-0.960
C ₇	0.378	0.393	0.378
C ₈	-0.394	-0.342	-0.388
C ₉	0.377	0.393	0.383
W	-1.051	-1.119	-1.040