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Population analysis using the SCF density.

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Orbital symmetries:

Occupied (SG) (SG) (SG) (SG) (PI) (PI) (SG)  
Virtual (PI) (PI) (SG)

The electronic state is 1-SG.

Alpha occ. eigenvalues -- -20.49362 -11.11511 -1.57552 -0.72113 -0.64241  
Alpha occ. eigenvalues -- -0.64241 -0.45542  
Alpha virt. eigenvalues -- 0.37448 0.37448 1.29137

Molecular Orbital Coefficients

		1	2	3	4	5
		(SG)--0	(SG)--0	(SG)--0	(SG)--0	(PI)--0
EIGENVALUES	--	-20.49362	-11.11511	-1.57552	-0.72113	-0.64241
1	1 C 1S	0.00006	0.99337	-0.15181	0.14468	0.00000
2	2S	-0.01256	0.02643	0.20128	-0.38176	0.00000
3	2PX	0.00000	0.00000	0.00000	0.00000	0.44260
4	2PY	0.00000	0.00000	0.00000	0.00000	0.00000
5	2PZ	-0.00944	0.00779	0.16987	-0.08086	0.00000
6	2 0 1S	0.99387	-0.00128	-0.22438	-0.13065	0.00000
7	2S	0.03212	-0.00656	0.74774	0.64875	0.00000
8	2PX	0.00000	0.00000	0.00000	0.00000	0.76636
9	2PY	0.00000	0.00000	0.00000	0.00000	0.00000
10	2PZ	-0.00937	-0.00201	-0.28069	0.71941	0.00000
		6	7	8	9	10
		(PI)--0	(SG)--0	(PI)--V	(PI)--V	(SG)--V
EIGENVALUES	--	-0.64241	-0.45542	0.37448	0.37448	1.29137
1	1 C 1S	0.00000	-0.17592	0.00000	0.00000	0.10417
2	2S	0.00000	0.84545	0.00000	0.00000	-1.31368
3	2PX	0.00000	0.00000	0.95804	0.00000	0.00000
4	2PY	0.44260	0.00000	0.00000	0.95804	0.00000
5	2PZ	0.00000	-0.55848	0.00000	0.00000	-1.44252
6	2 0 1S	0.00000	0.01997	0.00000	0.00000	-0.13789
7	2S	0.00000	-0.04794	0.00000	0.00000	1.49810
8	2PX	0.00000	0.00000	-0.72556	0.00000	0.00000
9	2PY	0.76636	0.00000	0.00000	-0.72556	0.00000
10	2PZ	0.00000	0.35882	0.00000	0.00000	-1.06409