
Population analysis using the SCF density.

Orbital symmetries:

Occupied (SGG) (SGU) (SGG) (SGU) (PIU) (PIU) (SGG)
Virtual (PIG) (PIG) (SGU)

The electronic state is 1-SGG.

Alpha occ. eigenvalues -- -15.55930 -15.55446 -1.54015 -0.71004 -0.64611
Alpha occ. eigenvalues -- -0.64611 -0.56496
Alpha virt. eigenvalues -- 0.32936 0.32936 1.38555

Molecular Orbital Coefficients

			1	2	3	4	5
			(SGG)--0	(SGU)--0	(SGG)--0	(SGU)--0	(PIU)--0
EIGENVALUES	--		-15.55930	-15.55446	-1.54015	-0.71004	-0.64611
1	1	N 1S	0.70267	0.70305	-0.18432	-0.16727	0.00000
2		2S	0.01280	0.03075	0.45930	0.72597	0.00000
3		2PX	0.00000	0.00000	0.00000	0.00000	0.00000
4		2PY	0.00000	0.00000	0.00000	0.00000	0.61138
5		2PZ	-0.00341	-0.01165	-0.26803	0.29611	0.00000
6	2	N 1S	0.70267	-0.70305	-0.18432	0.16727	0.00000
7		2S	0.01280	-0.03075	0.45930	-0.72597	0.00000
8		2PX	0.00000	0.00000	0.00000	0.00000	0.00000
9		2PY	0.00000	0.00000	0.00000	0.00000	0.61138
10		2PZ	0.00341	-0.01165	0.26803	0.29611	0.00000
			6	7	8	9	10
			(PIU)--0	(SGG)--0	(PIG)--V	(PIG)--V	(SGU)--V
EIGENVALUES	--		-0.64611	-0.56496	0.32936	0.32936	1.38555
1	1	N 1S	0.00000	-0.06513	0.00000	0.00000	-0.12360
2		2S	0.00000	0.42806	0.00000	0.00000	1.58382
3		2PX	0.61138	0.00000	0.86886	0.00000	0.00000
4		2PY	0.00000	0.00000	0.00000	0.86886	0.00000
5		2PZ	0.00000	0.60545	0.00000	0.00000	-1.38444
6	2	N 1S	0.00000	-0.06513	0.00000	0.00000	0.12360
7		2S	0.00000	0.42806	0.00000	0.00000	-1.58382
8		2PX	0.61138	0.00000	-0.86886	0.00000	0.00000
9		2PY	0.00000	0.00000	0.00000	-0.86886	0.00000
10		2PZ	0.00000	-0.60545	0.00000	0.00000	-1.38444