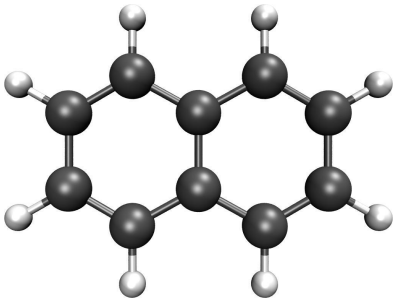
A Report On The Calculation Of The Single Point Energy Of Naphthalene At The PBE1PBE/6-31G(d,p) Level

osl - 15th June 2022



Abstract

The calculation of single point energy for the system 'Naphthalene' is presented, accompanied by automated analysis and image generation provided by the Silico software package. The calculation was performed using the Gaussian software package at the PBE1PBE/6-31G(d,p) level of theory. The total self-consistent field (SCF) energy of the system was found to be -10488.99 eV after 1 step. The highest-occupied molecular orbital (HOMO) and lowest-unoccupied molecular orbital (LUMO) were calculated to be -6.13 and -0.92 eV respectively, corresponding to a HOMO-LUMO band gap of 5.21 eV. The permanent dipole moment (PDM) was calculated to be 0.00 D.

Table 1: *Summary of overall calculation metadata. [a]: The date and time at which the calculation was completed. [b]: Total combined duration in real-time (wall-time) for all components of the calculation. [c]: Temperature used for thermochemistry analysis. [b]: Pressure used for thermochemistry analysis.*

Date[a]	Duration[b]	Success (Converged)	Computational package	Level of theory	Calculations	Wavefunction	Multiplicity	T[c] / K	p[d] / atm
15/06/2022 14:52:56	24 s	**True** (N/A)	Gaussian (2016+C.01)	PBE1PBE/ 6-31G(d,p)	Single Point	restricted	1 (singlet)	N/A	N/A

Summary Of Results

Scf Energy

Table 2: *Summary of SCF energy properties.*

No. of steps	1
Final energy	-10488.9903 eV
Final energy	-1,012,034 kJ·mol^{-1}

Geometry

Table 3: *Summary of geometry properties.*

Formula	$C_{10}H_8$
Exact mass	128.0626 g·mol^{-1}
Molar mass	128.1705 g·mol^{-1}
Alignment method	Minimal
X extension	6.74 Å
Y extension	4.97 Å
Z extension	0.00 Å
Linearity ratio	0.26
Planarity ratio	1.00

Molecular Orbitals

Table 4: *Summary of HOMO & LUMO properties.*

$\mathbf{E_{HOMO,LUMO}}$	5.21 eV
$\mathbf{E_{HOMO}}$	-6.13 eV
$\mathbf{E_{LUMO}}$	-0.92 eV

Permanent Dipole Moment

Table 5: *Summary of the permanent dipole moment properties.*

Total	0.00 D
X axis angle	0.00 °
XY plane angle	0.00 °

Methodology

Metadata

The calculation of the single point energy was performed using the **Gaussian (2016+C.01)** program, the **DFT** method with the **PBE1PBE** functional and the **6-31G(d,p)** basis set. It was completed on the **15th June 2022** after a total duration of **24 s** and **finished successfully**. The base multiplicity of the system under study was **1 (singlet)**. Finally, a **restricted wavefunction** was used, resulting in a single set of doubly occupied orbitals. The full calculation metadata is tabulated in table 1.

Analysis

The report presented here was generated using the Silico software package. This toolset relies upon a number of third-party applications and libraries which should be cited appropriately in derivative works. In particular, the calculation results described within were parsed by the cclib library.[1] Scientific constants which were used, among other things, for the interconversion of scientific units were provided by SciPy.[2] Three-dimensional plots of atom positions and calculated densities, including molecular orbitals, were rendered using Visual Molecular Dynamics (VMD)[3] and the Tachyon ray-tracer.[4] Finally, two-dimensional graphs were plotted using the MatPlotlib library,[5] while this report itself was prepared using the Mako template library[6] and the Weasyprint library[7], the latter of which was responsible for generation of the PDF file.

Discussion

Total SCF Energy

The total energy of the system was calculated at the **self-consistent field (SCF)** level, corresponding to the energy calculated by the density-functional theory (DFT) method, with a value of -10488.99 eV, corresponding to -1,012,034 KJmol^{-1}. A plot of the total SCF electron density is shown in figure 1.

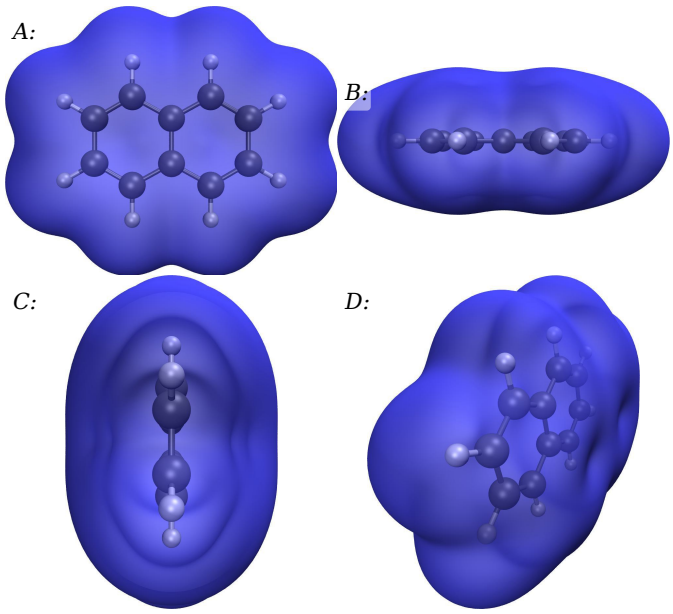


Figure 1: *Plot of the total SCF electron density, plotted with an isovalue of 0.0004. A: In the X/Y plane, B: In the X/Z plane, C: In the Z/Y plane, D: 45° to the axes.*

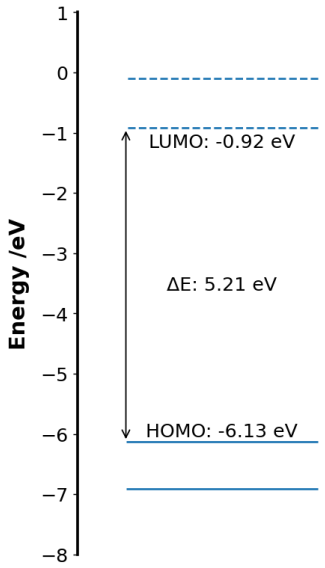


Figure 5: *Graph of the calculated molecular orbital energies in close proximity to the HOMO-LUMO gap. Solid lines: occupied orbitals, dashed lines: virtual orbitals.*

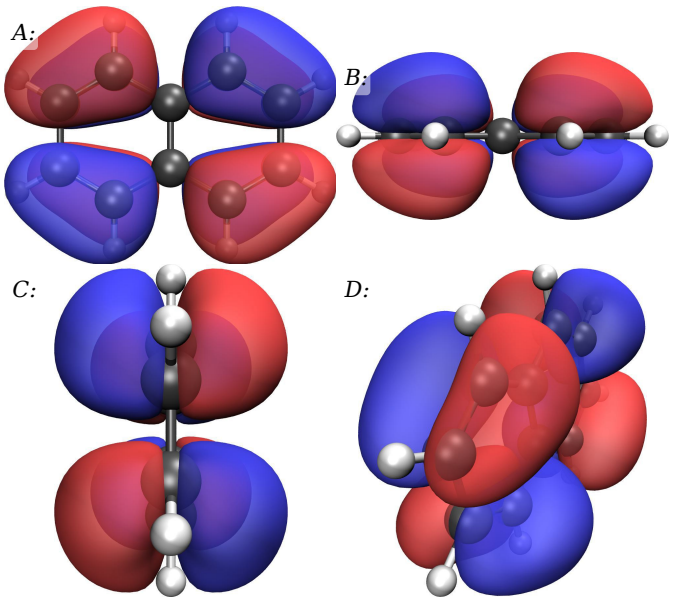


Figure 2: *Orbital density plots of the HOMO, plotted with isovalue: 0.02. A: In the X/Y plane, B: In the X/Z plane, C: In the Z/Y plane, D: 45° to the axes.*

Geometry

The **empirical formula** of the studied system was $C_{10}H_8$, corresponding to a **molecular mass** of 128.17 $gmol^{-1}$ and an **exact mass**, considering only specific atomic isotopes, of 128.06 $gmol^{-1}$. The molecular geometry was aligned to the cartesian (X, Y and Z) axes by the **Minimal (MIN)** method. Using this method, the **extent of the molecular system** in the X, Y and Z axes (L_X, L_Y and L_Z, corresponding to the molecular width, length and height respectively) was determined to be 6.74, 4.97 and 0.00 Å respectively. These extensions give rise to a **molecular linearity ratio** ($1-(L_Y/L_X)$) and **planarity ratio** ($1-(L_X/L_Y)$) of 0.26 and 1.00 respectively.

Permanent Dipole Moment

The calculated **permanent dipole moment** was exactly 0 D.

Molecular Orbitals

In total, 190 doubly occupied molecular orbitals were calculated, divided into 34 occupied orbitals and 156 unoccupied (or virtual) orbitals. The calculated energies of the **HOMO and LUMO** were -6.13 and -0.92 eV respectively, corresponding to a **HOMO-LUMO band gap** of 5.21 eV (figure 5). Plots of the orbital density for the HOMO and LUMO are shown in figures 2-3 respectively, while the orbital overlap between the HOMO and LUMO is shown in figure 4.

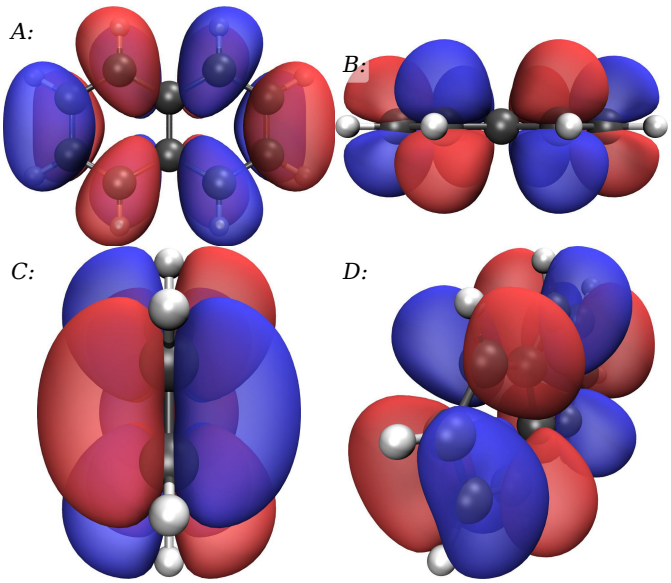


Figure 3: *Orbital density plots of the LUMO, plotted with isovalue: 0.02. A: In the X/Y plane, B: In the X/Z plane, C: In the Z/Y plane, D: 45° to the axes.*

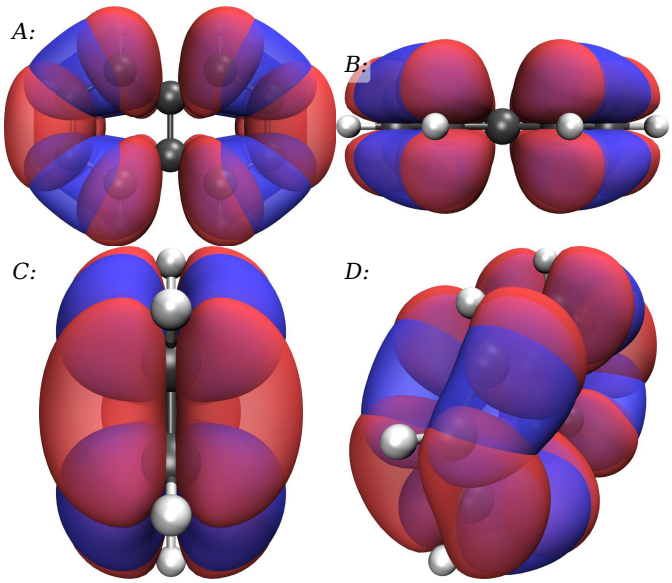


Figure 4: *Orbital density plots of the HOMO (blue) and LUMO (red), plotted simultaneously with isovalue: 0.02. A: In the X/Y plane, B: In the X/Z plane, C: In the Z/Y plane, D: 45° to the axes.*

Tables Of Results

Atom Coordinates

Table 6: *Coordinates of the atoms of the system under study, as aligned to the cartesian axes by the Minimal method.*

Element	X Coord /Å	Y Coord /Å	Z Coord /Å
C	-1.2404600	-1.3991400	0.0000000
C	-2.4260000	-0.7066400	0.0000000
C	-2.4260000	0.7066400	0.0000000
C	-1.2404600	1.3991400	-0.0000000
C	0.0000000	0.7142300	-0.0000000
C	-0.0000000	-0.7142300	0.0000000
C	1.2404600	-1.3991400	0.0000000
C	1.2404600	1.3991400	-0.0000000
C	2.4260000	0.7066400	-0.0000000
C	2.4260000	-0.7066400	-0.0000000
H	-1.2367000	-2.4862000	0.0000000
H	-3.3697000	-1.2439700	0.0000000
H	-3.3697000	1.2439700	0.0000000
H	-1.2367000	2.4862000	-0.0000000
H	1.2367000	-2.4862000	0.0000000
H	1.2367000	2.4862000	-0.0000000
H	3.3697000	1.2439700	-0.0000000
H	3.3697000	-1.2439700	-0.0000000

Molecular Orbitals

Table 7: *Energies of the calculated molecular orbitals.*

Level	Label	Symmetry	Energy /eV
50	LUMO+15	B1u	8.6396
49	LUMO+14	Ag	7.9114
48	LUMO+13	B1u	6.9150
47	LUMO+12	B3g	6.0692
46	LUMO+11	B2g	5.7949
45	LUMO+10	B2u	5.3487
44	LUMO+9	B1u	5.1506
43	LUMO+8	B3g	5.0031
42	LUMO+7	Ag	4.9519
41	LUMO+6	B1u	3.6912
40	LUMO+5	B2u	3.4207
39	LUMO+4	Au	2.9674
38	LUMO+3	Ag	2.9127
37	LUMO+2	B3u	1.0612
36	LUMO+1	B2g	-0.1010
35	**LUMO**	**B1g**	**-0.9244**
34	**HOMO**	**Au**	**-6.1307**
33	HOMO-1	B3u	-6.9084
32	HOMO-2	B2g	-8.0747
31	HOMO-3	Ag	-9.1879
30	HOMO-4	B1g	-9.2562
29	HOMO-5	B3g	-9.4032
28	HOMO-6	B2u	-10.2679
27	HOMO-7	B3u	-11.0274
26	HOMO-8	B1u	-11.1363
25	HOMO-9	B2u	-11.5961
24	HOMO-10	B3g	-11.6187
23	HOMO-11	Ag	-12.3015
22	HOMO-12	B1u	-12.4753
21	HOMO-13	Ag	-13.7777
20	HOMO-14	B3g	-14.2411
19	HOMO-15	B2u	-14.3709

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