# $C_{72}$: Gaudiene a Hollow and Aromatic All-Carbon Molecule

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## Cartesian coordinates

72
Energy =

| | | | |
|---|---|---|---|
| C | 0.0000000 | 2.7282389 | -3.7318211 |
| C | 1.1789225 | 2.0412217 | -4.0549321 |
| C | -1.1789225 | 2.0412217 | -4.0549321 |
| C | 2.7282389 | 0.0000000 | -3.7318211 |
| C | 2.0412217 | 1.1789225 | -4.0549321 |
| C | 2.0412217 | -1.1789225 | -4.0549321 |
| C | 1.1789225 | -2.0412217 | -4.0549321 |
| C | 0.0000000 | -2.7282389 | -3.7318211 |
| C | -1.1789225 | -2.0412217 | -4.0549321 |
| C | -2.0412217 | 1.1789225 | -4.0549321 |
| C | -2.7282389 | 0.0000000 | -3.7318211 |
| C | -2.0412217 | -1.1789225 | -4.0549321 |
| C | 0.0000000 | 3.7318211 | -2.7282389 |
| C | 3.7318211 | 0.0000000 | -2.7282389 |
| C | 0.0000000 | -3.7318211 | -2.7282389 |
| C | -3.7318211 | 0.0000000 | -2.7282389 |
| C | -4.0549321 | -1.1789225 | -2.0412217 |
| C | -4.0549321 | 1.1789225 | -2.0412217 |
| C | -1.1789225 | -4.0549321 | -2.0412217 |
| C | 1.1789225 | -4.0549321 | -2.0412217 |
| C | 4.0549321 | -1.1789225 | -2.0412217 |
| C | 4.0549321 | 1.1789225 | -2.0412217 |
| C | 1.1789225 | 4.0549321 | -2.0412217 |
| C | -1.1789225 | 4.0549321 | -2.0412217 |
| C | -2.0412217 | 4.0549321 | -1.1789225 |
| C | -4.0549321 | 2.0412217 | -1.1789225 |
| C | -4.0549321 | -2.0412217 | -1.1789225 |
| C | -2.0412217 | -4.0549321 | -1.1789225 |
| C | 2.0412217 | -4.0549321 | -1.1789225 |
| C | 4.0549321 | -2.0412217 | -1.1789225 |
| C | 4.0549321 | 2.0412217 | -1.1789225 |
| C | 2.0412217 | 4.0549321 | -1.1789225 |
| C | 2.7282389 | 3.7318211 | 0.0000000 |
| C | 3.7318211 | 2.7282389 | 0.0000000 |
| C | 3.7318211 | -2.7282389 | 0.0000000 |
| C | 2.7282389 | -3.7318211 | 0.0000000 |
| C | -2.7282389 | -3.7318211 | 0.0000000 |
| C | -3.7318211 | -2.7282389 | 0.0000000 |
| C | -3.7318211 | 2.7282389 | 0.0000000 |
| C | -2.7282389 | 3.7318211 | 0.0000000 |
| C | -2.7282389 | 0.0000000 | 3.7318211 |
| C | -2.0412217 | -1.1789225 | 4.0549321 |
| C | -2.0412217 | 1.1789225 | 4.0549321 |
| C | -1.1789225 | 2.0412217 | 4.0549321 |
| C | 0.0000000 | 2.7282389 | 3.7318211 |
| C | 1.1789225 | 2.0412217 | 4.0549321 |
| C | -1.1789225 | -2.0412217 | 4.0549321 |
| C | 0.0000000 | -2.7282389 | 3.7318211 |
| C | 1.1789225 | -2.0412217 | 4.0549321 |
| C | 2.0412217 | 1.1789225 | 4.0549321 |
| C | 2.7282389 | 0.0000000 | 3.7318211 |
| C | 2.0412217 | -1.1789225 | 4.0549321 |
| C | -3.7318211 | 0.0000000 | 2.7282389 |
| C | 0.0000000 | -3.7318211 | 2.7282389 |
| C | 0.0000000 | 3.7318211 | 2.7282389 |
| C | 3.7318211 | 0.0000000 | 2.7282389 |
| C | -1.1789225 | 4.0549321 | 2.0412217 |
| C | -4.0549321 | 1.1789225 | 2.0412217 |
| C | 1.1789225 | 4.0549321 | 2.0412217 |
| C | 4.0549321 | 1.1789225 | 2.0412217 |
| C | 4.0549321 | -1.1789225 | 2.0412217 |
| C | -1.1789225 | -4.0549321 | 2.0412217 |
| C | 1.1789225 | -4.0549321 | 2.0412217 |
| C | -4.0549321 | -1.1789225 | 2.0412217 |
| C | -4.0549321 | -2.0412217 | 1.1789225 |
| C | -2.0412217 | -4.0549321 | 1.1789225 |
| C | 2.0412217 | -4.0549321 | 1.1789225 |
| C | 4.0549321 | -2.0412217 | 1.1789225 |
| C | 4.0549321 | 2.0412217 | 1.1789225 |
| C | 2.0412217 | 4.0549321 | 1.1789225 |
| C | -2.0412217 | 4.0549321 | 1.1789225 |
| C | -4.0549321 | 2.0412217 | 1.1789225 |

Cartesian coordinates of $C_{72}$ in Å

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