

# Benzo[ghi]cyclopenta[cd]perylene

|                      |                                                                                   |
|----------------------|-----------------------------------------------------------------------------------|
| Inchi:               | InChI=1S/C24H12/c1-2-13-4-5-15-7-9-17-12-16-8-6-14-10-11-19-18(3-1)21(13)22(15)23 |
| InchiKey:            | MUDPUADNICLBBZ-UHFFFAOYSA-N                                                       |
| Formula:             | C24H12                                                                            |
| SMILES:              | C1=Cc2cc3ccc4ccc5cccc6c7ccc1c2c7c3c4c56                                           |
| Mol. weight [g/mol]: | 300.35                                                                            |
| CAS:                 | 190-88-5                                                                          |

## Physical Properties

| Property code | Value   | Unit    | Source         |
|---------------|---------|---------|----------------|
| gf            | 838.08  | kJ/mol  | Joback Method  |
| hf            | 650.53  | kJ/mol  | Joback Method  |
| hfus          | 41.06   | kJ/mol  | Joback Method  |
| hvap          | 82.54   | kJ/mol  | Joback Method  |
| log10ws       | -10.18  |         | Crippen Method |
| logp          | 6.812   |         | Crippen Method |
| mcvol         | 221.400 | ml/mol  | McGowan Method |
| pc            | 2370.28 | kPa     | Joback Method  |
| tb            | 890.88  | K       | Joback Method  |
| tc            | 1154.51 | K       | Joback Method  |
| tf            | 664.30  | K       | Joback Method  |
| vc            | 0.893   | m3/kmol | Joback Method  |

## Temperature Dependent Properties

| Property code | Value     | Unit    | Temperature [K] | Source        |
|---------------|-----------|---------|-----------------|---------------|
| cpg           | 625.52    | J/mol×K | 890.88          | Joback Method |
| cpg           | 639.68    | J/mol×K | 934.82          | Joback Method |
| cpg           | 654.66    | J/mol×K | 978.76          | Joback Method |
| cpg           | 670.86    | J/mol×K | 1022.70         | Joback Method |
| cpg           | 688.72    | J/mol×K | 1066.64         | Joback Method |
| cpg           | 708.65    | J/mol×K | 1110.57         | Joback Method |
| cpg           | 731.08    | J/mol×K | 1154.51         | Joback Method |
| dvisc         | 0.0266312 | Pa×s    | 664.30          | Joback Method |
| dvisc         | 0.0275768 | Pa×s    | 702.06          | Joback Method |

|       |           |      |        |               |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0284544 | Pa×s | 739.83 | Joback Method |
| dvisc | 0.0292708 | Pa×s | 777.59 | Joback Method |
| dvisc | 0.0300317 | Pa×s | 815.35 | Joback Method |
| dvisc | 0.0307426 | Pa×s | 853.12 | Joback Method |
| dvisc | 0.0314079 | Pa×s | 890.88 | Joback Method |

## Sources

|                        |                                                                                                                                           |
|------------------------|-------------------------------------------------------------------------------------------------------------------------------------------|
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>                     |
| <b>NIST Webbook:</b>   | <a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C190885&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C190885&amp;Units=SI</a> |
| <b>Crippen Method:</b> | <a href="http://pubs.acs.org/doi/abs/10.1021/ci990307l">http://pubs.acs.org/doi/abs/10.1021/ci990307l</a>                                 |
| <b>Crippen Method:</b> | <a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>                         |
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                                     |

## Legend

|                 |                                                 |
|-----------------|-------------------------------------------------|
| <b>cpg:</b>     | Ideal gas heat capacity                         |
| <b>dvisc:</b>   | Dynamic viscosity                               |
| <b>gf:</b>      | Standard Gibbs free energy of formation         |
| <b>hf:</b>      | Enthalpy of formation at standard conditions    |
| <b>hfus:</b>    | Enthalpy of fusion at standard conditions       |
| <b>hvap:</b>    | Enthalpy of vaporization at standard conditions |
| <b>log10ws:</b> | Log10 of Water solubility in mol/l              |
| <b>logp:</b>    | Octanol/Water partition coefficient             |
| <b>mcvol:</b>   | McGowan's characteristic volume                 |
| <b>pc:</b>      | Critical Pressure                               |
| <b>tb:</b>      | Normal Boiling Point Temperature                |
| <b>tc:</b>      | Critical Temperature                            |
| <b>tf:</b>      | Normal melting (fusion) point                   |
| <b>vc:</b>      | Critical Volume                                 |

Latest version available from:

<https://www.chemeo.com/cid/52-890-6/Benzo-ghi-cyclopenta-cd-perylene.pdf>

Generated by Cheméo on 2025-12-24 00:46:33.785411757 +0000 UTC m=+6285391.315452411.

Cheméo (<https://www.chemeo.com>) is the biggest free database of chemical and physical data for the process industry.