

Ergost-5-en-3-ol, acetate, (3«beta»,24R)-

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|-----------------------------|--|
| Other names: | Campesterol acetate |
| Inchi: | InChI=1S/C30H50O2/c1-19(2)20(3)8-9-21(4)26-12-13-27-25-11-10-23-18-24(32-22(5)31 |
| InchiKey: | JOBAYBRAHVTSSW-VYOIUQHLSA-N |
| Formula: | C30H50O2 |
| SMILES: | CC(=O)OC1CCC2(C)C(=CCC3C2CCC2(C)C(C(C)CCC(C)C(C)C)CCC32)C1 |
| Mol. weight [g/mol]: | 442.72 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|---------|----------------|
| gf | 129.20 | kJ/mol | Joback Method |
| hf | -647.00 | kJ/mol | Joback Method |
| hfus | 39.16 | kJ/mol | Joback Method |
| hvap | 88.60 | kJ/mol | Joback Method |
| log10ws | -8.61 | | Crippen Method |
| logp | 8.206 | | Crippen Method |
| mcvol | 393.260 | ml/mol | McGowan Method |
| pc | 880.00 | kPa | Joback Method |
| rinpol | 3323.00 | | NIST Webbook |
| rinpol | 3278.00 | | NIST Webbook |
| tb | 999.69 | K | Joback Method |
| tc | 1231.31 | K | Joback Method |
| tf | 557.54 | K | Joback Method |
| vc | 1.488 | m3/kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|---------|---------|-----------------|---------------|
| cpg | 1504.06 | J/molxK | 999.69 | Joback Method |
| cpg | 1537.82 | J/molxK | 1038.29 | Joback Method |
| cpg | 1572.22 | J/molxK | 1076.90 | Joback Method |
| cpg | 1607.63 | J/molxK | 1115.50 | Joback Method |
| cpg | 1644.38 | J/molxK | 1154.10 | Joback Method |
| cpg | 1682.84 | J/molxK | 1192.71 | Joback Method |
| cpg | 1723.35 | J/molxK | 1231.31 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U141997&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci990307l |
| Crippen Method: | https://www.cheméo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |

Legend

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

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