

# Selina-4,11-dien-14-al

|                      |                                                                                      |
|----------------------|--------------------------------------------------------------------------------------|
| Inchi:               | InChI=1S/C15H22O/c1-11(2)12-6-8-15(3)7-4-5-13(10-16)14(15)9-12/h10,12H,1,4-9H2,2H,3H |
| InchiKey:            | VNYQZKHAWVXJCR-WFASDCNBSA-N                                                          |
| Formula:             | C15H22O                                                                              |
| SMILES:              | C=C(C)C1CCC2(C)CCCC(C=O)=C2C1                                                        |
| Mol. weight [g/mol]: | 218.33                                                                               |

## Physical Properties

| Property code | Value   | Unit   | Source         |
|---------------|---------|--------|----------------|
| hfus          | 16.32   | kJ/mol | Joback Method  |
| logp          | 4.048   |        | Crippen Method |
| mcvol         | 193.460 | ml/mol | McGowan Method |
| rinpol        | 1758.00 |        | NIST Webbook   |
| rinpol        | 1758.00 |        | NIST Webbook   |
| rinpol        | 1758.00 |        | NIST Webbook   |

## Temperature Dependent Properties

| Property code | Value  | Unit    | Temperature [K] | Source        |
|---------------|--------|---------|-----------------|---------------|
| cpg           | 526.25 | J/mol×K | 627.74          | Joback Method |
| cpg           | 546.54 | J/mol×K | 665.73          | Joback Method |
| cpg           | 565.61 | J/mol×K | 703.71          | Joback Method |
| cpg           | 583.64 | J/mol×K | 741.70          | Joback Method |
| cpg           | 600.80 | J/mol×K | 779.69          | Joback Method |
| cpg           | 617.25 | J/mol×K | 817.67          | Joback Method |
| cpg           | 633.16 | J/mol×K | 855.66          | Joback Method |

## Sources

Cheméo Relay (1.0):

Cheméo Relay (1.0, beta):

NIST Webbook:

<http://webbook.nist.gov/cgi/cbook.cgi?ID=R73335&Units=SI>

|                        |                                                                                                                       |
|------------------------|-----------------------------------------------------------------------------------------------------------------------|
| <b>Joback Method:</b>  | <a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>                 |
| <b>McGowan Method:</b> | <a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</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>     |

## Legend

|                |                                           |
|----------------|-------------------------------------------|
| <b>cpg:</b>    | Ideal gas heat capacity                   |
| <b>hfus:</b>   | Enthalpy of fusion at standard conditions |
| <b>logp:</b>   | Octanol/Water partition coefficient       |
| <b>mcvol:</b>  | McGowan's characteristic volume           |
| <b>rinpol:</b> | Non-polar retention indices               |

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