

# «alpha»-D-Galactopyranose, 1,2:3,4-bis-O-(1-methylethylidene)-

Other names:

Galactopyranose, 1,2:3,4-di-O-isopropylidene-, «alpha»-D-

Diisopropylidenegalactose

1,2:3,4-di-O-Isopropylidene-D-galactopyranose

1,2,3,4-di-O-Isopropylidene-D-galactopyranose

1,2,3,4-di-O-Isopropylidenegalactopyranose

Galactopyranose, 1,2:3,4-di-o-isopropylidene-, d-

Inchi: InChI=1S/C12H20O6/c1-11(2)15-7-6(5-13)14-10-9(8(7)16-11)17-12(3,4)18-10/h6-10,13H

InchiKey: POORJMIIHXHXAV-MBXMOIHESA-N

Formula: C12H20O6

SMILES: CC1(C)OC2OC(CO)C3OC(C)(C)OC3C2O1

Mol. weight [g/mol]: 260.28

CAS: 4064-06-6

## Physical Properties

Property code	Value	Unit	Source
gf	-413.13	kJ/mol	Joback Method
hf	-954.20	kJ/mol	Joback Method
hfus	50.61	kJ/mol	Joback Method
hvap	78.25	kJ/mol	Joback Method
log10ws	-1.50		Crippen Method
logp	0.375		Crippen Method
mcvol	182.580	ml/mol	McGowan Method
pc	2770.08	kPa	Joback Method
tb	715.72	K	Joback Method
tc	928.65	K	Joback Method
tf	492.77	K	Joback Method
vc	0.670	m <sup>3</sup> /kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	607.20	J/mol×K	715.72	Joback Method
cpg	623.91	J/mol×K	751.21	Joback Method
cpg	640.18	J/mol×K	786.70	Joback Method

cpg	656.22	J/mol×K	822.19	Joback Method
cpg	672.24	J/mol×K	857.67	Joback Method
cpg	688.45	J/mol×K	893.16	Joback Method
cpg	705.06	J/mol×K	928.65	Joback Method

## Sources

<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>
<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=C4064066&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C4064066&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<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/93-752-4/alpha-D-Galactopyranose-1-2-3-4-bis-O-1-methylethylidene.pdf>

Generated by Cheméo on 2024-04-25 19:07:35.823742522 +0000 UTC m=+16361304.744319844.

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