

Other names

undecaethylene glycol, bis(neptafluorobutylate)

Property code	Value	Unit	Temperature [K]	Source
cpg	1941.76	J/mol×K	1232.98	Joback Method
cpg	1941.05	J/mol×K	1342.87	Joback Method
cpg	1921.64	J/mol×K	1452.77	Joback Method
cpg	1884.95	J/mol×K	1562.66	Joback Method
cpg	1832.43	J/mol×K	1672.55	Joback Method
cpg	1765.51	J/mol×K	1782.45	Joback Method
cpg	1685.61	J/mol×K	1892.34	Joback Method

Sources

Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U352033&Units=SI
Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws

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

Latest version available from:

<https://www.chemeo.com/cid/44-609-7/2-2-2-2-2-2-2-2-2-2-2-3-3-4-4-4-Heptafluorobutanoyl-oxyethoxy-ethoxy-et>

Generated by Cheméo on 2025-12-06 04:22:18.717547158 +0000 UTC m=+4743136.247587821.

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