

### Surface Energy Data for PVF: Poly(vinyl fluoride), CAS # 24981-14-4

Source <sup>(a)</sup>	Mst. Type <sup>(b)</sup>	Data <sup>(c)</sup>	Comments <sup>(d)</sup>
Ellison, 1954 <sup>(8)</sup>	Critical ST	$\gamma_c = 28 \text{ mJ/m}^2$ , 20°C	Various test liquids.
Wu, 1971 <sup>(29)</sup>	Contact angle	$\theta_W^Y = 80^\circ$ , 20°C	
Moshonov, 1980 <sup>(118)</sup>	Contact angle	$\theta_W^Y = 85^\circ$ ; no temp cited	Measured 60 secs. after application of water droplet; surface cleaned with isopropanol at 60°C and rinsed with methanol.
Vargha-Butler, 1985 <sup>(180)</sup>	Contact angle	$\theta_W^A = 88.6^\circ$ ; 20°C	
Wu, 1971 <sup>(29)</sup>	Contact angle	$\gamma_s = 36.7 \text{ mJ/m}^2$ ( $\gamma_s^d = 32.0$ , $\gamma_s^p = 4.7$ ); 20°C	Test liquids: water and diiodomethane, by geometric mean equation.
Wu, 1971 <sup>(29)</sup>	Contact angle	$\gamma_s = 38.4 \text{ mJ/m}^2$ ( $\gamma_s^d = 27.3$ , $\gamma_s^p = 11.1$ ); 20°C	Test liquids: water and diiodomethane, by harmonic mean equation.
Kitazaki, 1972 <sup>(191)</sup>	Contact angle	$\gamma_s = 43.5 \text{ mJ/m}^2$ ( $\gamma_s^d = 42.3$ , $\gamma_s^p = 1.2$ ); no temp cited	Various test liquids; original results split polar component into hydrogen- and non-hydrogen bonding parameters.
Wu, 1979 <sup>(45)</sup>	Contact angle	$\gamma_c = 37.5 \text{ mJ/m}^2$ , 20°C	Test liquids not known; calculated by the equation of state method.
Brewis, 1981 <sup>(261)</sup>	Contact angle	$\gamma_s = 36.7 \text{ mJ/m}^2$ ( $\gamma_s^d = 36.7$ , $\gamma_s^p = 0.0$ ); no temp cited	Test liquids not known.
Vargha-Butler, 1985 <sup>(180)</sup>	Contact angle	$\gamma_c = 29.4 \text{ mJ/m}^2$ ; 20°C	Test liquids not known; calculated by the equation of state method.
Lloyd, 1995 <sup>(218)</sup>	Contact angle	$\gamma_s = 36.7 \text{ mJ/m}^2$ ( $\gamma_s^{LW} = 34.8$ , $\gamma_s^{AB} = 1.9$ , $\gamma_s^+ = 0.2$ , $\gamma_s^- = 4.5$ ); no temp cited	Test liquids not known; acid-base analysis. Tedlar.
Lloyd, 1995 <sup>(218)</sup>	Contact angle	$\gamma_s = 39.0 \text{ mJ/m}^2$ ( $\gamma_s^{LW} = 38.0$ , $\gamma_s^{AB} = 1.0$ , $\gamma_s^+ = 0.02$ , $\gamma_s^- = 13.2$ ); no temp cited	Test liquids not known; acid-base analysis.
Morra, 1999 <sup>(134)</sup>	Contact angle	$\gamma_s = 36.8 \text{ mJ/m}^2$ ( $\gamma_s^{LW} = 34.1$ , $\gamma_s^{AB} = 2.7$ , $\gamma_s^+ = 8.3$ , $\gamma_s^- = 0.2$ ); no temp cited	Test liquids not known; acid-base analysis based on reference values for water of $\gamma^+ = 48.5 \text{ mJ/m}^2$ and $\gamma^- = 11.2 \text{ mJ/m}^2$ .
Uschold, 1999 <sup>(211)</sup>	Contact angle	$\gamma_s = 38 \text{ mJ/m}^2$ ; no temp cited	
Chang, 2000 <sup>(162)</sup>	Contact angle	$\gamma_s = 35.2 \text{ mJ/m}^2$ ; no temp cited	
Kwok, 2000 <sup>(166)</sup>	Contact angle	$\gamma_c = 35.7 \text{ mJ/m}^2$ ; no temp cited	Re-calculated by equation of state method from data produced by Ellison, 1954 <sup>(8)</sup> .
Wu, 1971 <sup>(29)</sup>	From polymer melt	$\gamma_s = 37.5 \text{ mJ/m}^2$ ( $\gamma_s^d = 31.6$ , $\gamma_s^p = 5.9$ ); 20°C	Direct measurement of polymer melt extrapolated to 20°C.
Lee, 1968 <sup>(131)</sup>	Calculated	$\gamma_s = 41 \text{ mJ/m}^2$ ; no temp cited	Calculated from glass temperature of 323K.
Wu, 1968 <sup>(182)</sup>	Calculated	$\gamma_s = 29 \text{ mJ/m}^2$ ; 20°C	Calculated from molecular constitution.
Van Krevelen, 1976 <sup>(85)</sup>	Calculated	$\gamma_s = 32.5 \text{ mJ/m}^2$ ; no temp cited	Calculated from parachor parameter.
Vargha-Butler, 1985 <sup>(180)</sup>	Calculated	$\gamma_s = 28.8 \text{ mJ/m}^2$ ; no temp cited	Calculated from sedimentation volume.