

Supplemental Information for Reference Correlation of the Viscosity of Argon

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Additional Details on the Fitting Procedure

The program Eureka¹ was used with the error metric “absolute error” selected. The formula building blocks were constant, input variable, addition, subtraction, multiplication and division. No variable preparation options were selected. The data were entered as T_r , ρ_r , and the function $F(\rho_r, T_r)$ as described in the text. The function F was set to zero for all points with density less than 2 mol/l in order to force the function to reproduce the theoretical behavior at low density. Due to this, many points at low density in the primary set were not used in the fitting process. Weights were obtained by a trial-and-error process with the goal of representing the majority of the data to within or near the estimated experimental uncertainty. Accordingly, a subset of the primary data listed in Table 2 in the main manuscript were assigned the following weights:

1st author	wt	Notes	1st author	wt	Notes
Zhou	5	vapor points	Mostert	5	
Zhou	8	liquid points	Diller	8	
Humberg	1		Haynes	10	
Berg	1		Vermesse	3	
Abramson	0.5		Rabinovich	10	
Vogel	20		Gracki	1	
Xiao	20		Boon	7	
Hurly	1		Flynn	7	
Evers	3		Flynn	10	Points with $T_r > 1.49$ and $\rho_r > 0.8$
Wilhelm	10				

In addition, to guide the extrapolation behavior at 2000 K, the following points, estimated from a corresponding states method, were added to the fitting procedure.

T_r	ρ_r	F	wt	T_r	ρ_r	F	wt
13.2725	0.6965	3.429	50	13.2725	1.8469	8.756	50
13.2725	0.9435	4.202	50	13.2725	1.9481	9.511	50
13.2725	1.1495	4.968	50	13.2725	2.6656	17.273	50
13.2725	1.3255	5.371	50	13.2725	3.1301	25.750	50
13.2725	1.4789	6.491	50	13.2725	3.4813	35.050	50
13.2725	1.6147	7.248	50	13.2725	3.7655	45.109	50
13.2725	1.7365	8.002	50				

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