

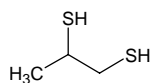
# 1,2-propanedithiol

## 1,2-プロパンジチオール

Formula C<sub>3</sub>H<sub>8</sub>S<sub>2</sub>

Mol. Wt. 108.23

CAS No. 814-67-5



SEQ No. 2196 チオール類

FEMA No. 3520

JECFA No. 536

	JFFMA	JECFA	FCC
Chemical name	1,2-propanedithiol	1,2-Propanedithiol	
Content (NLT)	95.0 % (GC)	96 %	
Refractive index	1.526-1.538 (n <sub>20D</sub> )	1.531-1.541 (n <sub>20D</sub> )	
Specific gravity	1.058-1.068 (d <sub>20/20</sub> )	1.075-1.081 (d <sub>25/25</sub> )	
Acid value (NMT)			
Melting/Congeeing point			
Angular/Specific rotation			
Heavy metals (NMT)			
Solubility in ethanol			
Other requirements		Chlorinated monothiol: max. 1.1%	
Identification*	IR : 1,3,5 MS : 3,4,5 NMR : 1,3	IR, NMR	

### \*References

- 1: Joint FAO/WHO Expert Committee on Food Additive (JECFA)
- 2: Food Chemicals Codex (FCC)
- 3: 有機化合物のスペクトルデータベース SDBS(独立行政法人産業技術総合研究所)
- 4: Wiley's Registry of Mass spectral Database
- 5: NIST/EPA/NIH Mass Spectral Library
- 6: Sigma-Aldrich