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PYRROMETHENE 597-8C9

Synonym: 2,6-di-tert-butyl-8-nonyl-1,3,5,7-tetramethylpyrromethene- BF₂ Complex Catalog No.: 05971 CAS Registry Number: N/A Chemical Formula: C₃₀H₄₉N₂BF₂ MW: 486.54 Appearance: Red-orange solid Melting Point: 247°C



Suggested Use: lithography, fl hydrocarbon probe or taggant, electro-luminescence, DVD (digital video data)

Solubility Limits (g/l) (room temperature):

Ethanol 3.7 Heptane ≥10.3 Ethyl Acetate very soluble Acetonitrile very soluble p-Dioxane very soluble EPH PPH Toluene THF Dow Corning 200 Silicone Oil 0.47

very soluble very soluble very soluble very soluble

Spectral Information:

 $\lambda_{\text{max,abs}} = 524$ nm (ethanol), 525 nm (methylene chloride), 527.8 nm (diesel fuel) $\lambda_{\text{max,fl}} = 588$ nm (ethanol), 585 nm (methylene chloride), 590 nm (diesel fuel) $\epsilon_{524} = 5.25 \times 10^4$ liter mol⁻¹ cm⁻¹



The information presented above is believed to be accurate but is not a specification. The customer is fully responsible for determining the suitability of this product for use in their application. Exciton, Inc. does not represent that the information is sufficient or complete for any specific application.

Pyrromethene 597-8C9 in Methanol