

## Product Datasheet

### 5-Propargylamino-ddUTP (ATTO-Thio12) (orb64785)

|                          |                                                                                                                                                             |
|--------------------------|-------------------------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Description</b>       | 5-Propargylamino-ddUTP conjugated to ATTO-Thio12.                                                                                                           |
| <b>Form/Appearance</b>   | solution in water; pH: 7.5 ± 0.5                                                                                                                            |
| <b>Concentration</b>     | 0.50 mM-0.55 mM                                                                                                                                             |
| <b>Storage</b>           | store at -20 °C                                                                                                                                             |
| <b>Note</b>              | For research use only                                                                                                                                       |
| <b>Application notes</b> | <b>Spectroscopic Propertie:</b> λ <sub>exc</sub> 582 nm, λ <sub>em</sub> 607 nm, ε 110.0 L mmol <sup>-1</sup> cm <sup>-1</sup> (Tris-HCl pH 7.5).           |
| <b>Formula</b>           | C <sub>41</sub> H <sub>47</sub> N <sub>6</sub> O <sub>15</sub> P <sub>3</sub> S                                                                             |
| <b>Purity</b>            | ≥ 95% (HPLC)                                                                                                                                                |
| <b>MW</b>                | Theoretical MW: 988.83 g/mol (free acid)                                                                                                                    |
| <b>SMILES</b>            | <chem>C(C#Cc1cn(c(=O)[nH]c1=O)[C@@H]1O[C@H](COP(=O)(O)OP(=O)(O)OP(=O)(O)[O-])CC1)NC(=O)CCCN(C(=O)c1c(ccc1)c1c2c(sc3c1ccc(c3)N(C)C)cc(=[N+](C)C)cc2)C</chem> |
| <b>Expiration Date</b>   | 12 months from date of receipt.                                                                                                                             |