

## Product Datasheet

### 7-Propargylamino-7-deaza-ddGTP (ATTO-Rho11) (orb64599)

|                          |                                                                                                                                                                                  |
|--------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Description</b>       | 7-Propargylamino-7-deaza-ddGTP conjugated to ATTO-Rho11.                                                                                                                         |
| <b>Form/Appearance</b>   | solution in water; pH: 7.5 ± 0.5                                                                                                                                                 |
| <b>Concentration</b>     | 1.0 mM-1.1 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> 572 nm, λ <sub>em</sub> 595 nm, ε 120.0 L mmol <sup>-1</sup> cm <sup>-1</sup> (Tris-HCl pH 7.5).                                |
| <b>Formula</b>           | C <sub>49</sub> H <sub>57</sub> N <sub>8</sub> O <sub>16</sub> P <sub>3</sub>                                                                                                    |
| <b>Purity</b>            | ≥ 95% (HPLC)                                                                                                                                                                     |
| <b>MW</b>                | Theoretical MW: 1106.96 g/mol (free acid); Detected MW: 1106.31 g/mol (free acid)                                                                                                |
| <b>SMILES</b>            | <chem>N(CC#Cc1c2c(n(c1)[C@@H]1O[C@H](COP(=O)(O)OP(=O)(O)OP(=O)(O)O)[C@@H](O)C1)nc(N)[nH]c2=O)C(=O)CCCN(C(=O)c1c(C2=c3cc4c(=[N+](CCC4)CC)cc3Oc3c2cc2c(c3)N(CCC2)CC)cccc1)C</chem> |
| <b>Expiration Date</b>   | 12 months from date of receipt.                                                                                                                                                  |