

## Product Datasheet

### 5-Propargylamino-ddCTP (ATTO-390) (orb64860)

|                          |                                                                                                                                                         |
|--------------------------|---------------------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Description</b>       | 5-Propargylamino-ddCTP conjugated to ATTO-390.                                                                                                          |
| <b>Form/Appearance</b>   | solution in water; Color: slightly yellow; 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> $\lambda_{exc}$ 390 nm, $\lambda_{em}$ 476 nm, $\epsilon$ 24.0 L mmol <sup>-1</sup> cm <sup>-1</sup> (Tris-HCl pH 7.5). |
| <b>Formula</b>           | C <sub>32</sub> H <sub>42</sub> N <sub>5</sub> O <sub>15</sub> P <sub>3</sub>                                                                           |
| <b>Purity</b>            | ≥ 95% (HPLC)                                                                                                                                            |
| <b>MW</b>                | Theoretical MW: 829.63 g/mol (free acid); Detected MW: 829.19 g/mol (free acid)                                                                         |
| <b>SMILES</b>            | <chem>C1C[C@@H](O[C@@H]1COP(=O)(OP(=O)(OP(=O)(O)O)O)O)n1c(=O)nc(c(c1)C#CCNC(=O)CCCN1c2cc3oc(=O)cc(C)c3cc2C(C)CC1(C)C)N</chem>                           |
| <b>Expiration Date</b>   | 12 months from date of receipt.                                                                                                                         |