

## Product Datasheet

### 5-Propargylamino-ddUTP (ATTO-532) (orb64809)

|                          |                                                                                                                                                                                    |
|--------------------------|------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Description</b>       | 5-Propargylamino-ddUTP conjugated to ATTO-532.                                                                                                                                     |
| <b>Form/Appearance</b>   | solution in water; Color: orange; 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> 532 nm, λ <sub>em</sub> 552 nm, ε 115.0 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>22</sub> P <sub>3</sub> S <sub>2</sub>                                                                                       |
| <b>Purity</b>            | ≥ 95% (HPLC)                                                                                                                                                                       |
| <b>MW</b>                | Theoretical MW: 1132.89 g/mol (free acid); Detected MW: 1132.14 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)[nH]c(=O)c(c1)C#CCNC(=O)CCCN(C(=O)c1c(ccc1)c1c2c(oc3c1cc/c(=[NH+]/CC)/c3S(=O)(=O)[O-])c(c(cc2)NCC)S(=O)(=O)O)C</chem> |
| <b>Expiration Date</b>   | 12 months from date of receipt.                                                                                                                                                    |