

## Product Datasheet

### 5/6-Carboxyrhodamine 110-PEG3-Azide (orb1734905)

|                          |                                                                                                                                                                                  |
|--------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|
| <b>Description</b>       | With excitation maximum at 501 nm, the azide-carboxyrhodamine 110 conjugate is an excellent match to...                                                                          |
| <b>Form/Appearance</b>   | solid; Color: dark red                                                                                                                                                           |
| <b>Storage</b>           | store at -20 °C                                                                                                                                                                  |
| <b>Note</b>              | For research use only                                                                                                                                                            |
| <b>Application notes</b> | <b>Spectroscopic Propertie</b> : λabs 501 nm, λem 525 nm, ε 74.0 L mmol <sup>-1</sup> cm <sup>-1</sup> (in MeOH).                                                                |
| <b>Formula</b>           | C <sub>29</sub> H <sub>30</sub> N <sub>6</sub> O <sub>7</sub>                                                                                                                    |
| <b>Purity</b>            | ≥ 90% (HPLC)                                                                                                                                                                     |
| <b>MW</b>                | Theoretical MW: 574.59 g/mol; Detected MW: 574.21 g/mol                                                                                                                          |
| <b>SMILES</b>            | <chem>N(=[N+]=[N-])CCOCCOCCOCCNC(=O)c1cc(c2c3ccc(cc3oc3cc(=[NH2+])ccc23)N)c(cc1)C(=O)[O-].O=C(NCCOCCOCCOCCN=[N+]=[N-])c1cc(c2c3ccc(cc3oc3cc(=[NH2+])ccc23)N)cc1)C(=O)[O-]</chem> |
| <b>Solubility (25°C)</b> | DMF, DMSO, MeOH                                                                                                                                                                  |
| <b>Expiration Date</b>   | 12 months from date of receipt.                                                                                                                                                  |