



## Product Information

### Tri-GalNAc(OAc)3, Purity ≥98%

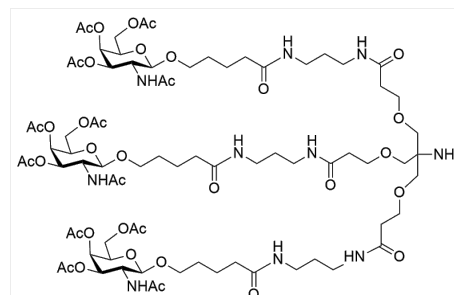
**Cat. No.:** X24-11-YM016

**Size:** 1 g; 2 g; 5 g; 10 g

**CAS Number:** 1159408-64-6

**Synonym:** 1159408-64-6; Peracetylated tri-GalNAc-amine

**This product is for research use only and is not intended for diagnostic use.**



#### Product Information

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| <b>Description</b>       | Tri-GalNAc(OAc)3 is used to synthesize conjugates for targeted drug delivery systems. These conjugates can specifically bind to receptors on target cells, such as the asialoglycoprotein receptor (ASGPR), ensuring the drugs reach their intended targets with high precision.                                                                                                                                                                                                                                                                                                                                                    |
| <b>Molecular Weight</b>  | 1793.9                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                              |
| <b>Molecular Formula</b> | C <sub>79</sub> H <sub>128</sub> N <sub>10</sub> O <sub>36</sub>                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    |
| <b>Targets</b>           | ASGPR                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                               |
| <b>IUPAC Name</b>        | [(2R,3R,4R,5R,6R)-5-Acetamido-6-[5-[3-[3-[3-[3-[3-[5-[(2R,3R,4R,5R,6R)-3-acetamido-4,5-diacetyloxy-6-(acetyloxymethyl)oxan-2-yl]oxypentanoylamino]propylamino]-3-oxopropoxy]-2-[[3-[3-[5-[(2R,3R,4R,5R,6R)-3-acetamido-4,5-diacetyloxy-6-(acetyloxymethyl)oxan-2-yl]oxypentanoylamino]propylamino]-3-oxopropoxy]methyl]-2-aminopropoxy]propanoylamino]propylamino]-5-oxopentoxo]-3,4-diacetyloxyoxan-2-yl]methyl acetate                                                                                                                                                                                                            |
| <b>InChI</b>             | InChI=1S/C79H128N10O36/c1-46(90)87-67-73(120-55(10)99)70(117-52(7)96)58(40-114-49(4)93)123-76(67)111-34-16-13-22-61(102)81-28-19-31-84-64(105)25-37-108-43-79(80,44-109-38-26-65(106)85-32-20-29-82-62(103)23-14-17-35-112-77-68(88-47(2)91)74(121-56(11)100)71(118-53(8)97)59(124-77)41-115-50(5)94)45-110-39-27-66(107)86-33-21-30-83-63(104)24-15-18-36-113-78-69(89-48(3)92)75(122-57(12)101)72(119-54(9)98)60(125-78)42-116-51(6)95/h58-60,67-78H,13-45,80H2,1-12H3,(H,81,102)(H,82,103)(H,83,104)(H,84,105)(H,85,106)(H,86,107)(H,87,90)(H,88,91)(H,89,92)/t58-,59-,60-,67-,68-,69-,70+,71+,72+,73-,74-,75-,76-,77-,78-/m1/s1 |
| <b>InChI Key</b>         | VGHVGBJYGYFWED-PNPDLPPQSA-N                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         |
| <b>SMILES string</b>     | CC(=O)N[C@@H]1[C@H]([C@H]([C@H]([C@H]([C@H]([C@H]1O)CCCCC(=O)NCCCNC(=O)CCOCC(COCCC(=O)NCCCNC(=O)CCCCO[C@H]2[C@@H]([C@H]([C@H]([C@H]([C@H]([C@H]2O)COC(=O)C)OC(=O)C)OC(=O)C)NC(=O)C)(COCCC(=O)NCCCNC(=O)CCCCO[C@H]3[C@@H]([C@H]([C@H]([C@H]([C@H]([C@H]3O)COC(=O)C)OC(=O)C)OC(=O)C)NC(=O)C)N)COC(=O)C)OC(=O)C)OC(=O)C                                                                                                                                                                                                                                                                                                                |
| <b>Form</b>              | Solid                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                               |



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| <b>Purity</b>        | ≥98%, determined by HPLC.                                                                      |
| <b>Impurities</b>    | Free from inappropriate visible particulates, foreign matter, discoloration, or other defects. |
| <b>Identity</b>      | Conforms to structure, determined by <sup>1</sup> H NMR.                                       |
| <b>Test Method</b>   | NMR; HPLC                                                                                      |
| <b>Quality Level</b> | Research grade                                                                                 |
| <b>Storage</b>       | Store under dry conditions at room temperature.                                                |

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