

## Supplementary Information for

### Low Valency Display of Glycomimetic Inhibitor of Adhesion of Pathogenic *Candida albicans* to Human Buccal Epithelial Cells

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#### Content

1.1. <sup>1</sup>H NMR of *N, N'*-di-(2,3,4,6-tetra-*O*-acetyl-β-*D*-galactopyranosyl-1,2,3-triazol-4-ylmethylamide)-*N''*-(2-bromoacetamido)-5-aminobenzene-1,3-dicarboxamide (**7**)

1.2. <sup>13</sup>C NMR of *N, N'*-di-(2,3,4,6-tetra-*O*-acetyl-β-*D*-galactopyranosyl-1,2,3-triazol-4-ylmethylamide)-*N''*-(2-bromoacetamido)-5-aminobenzene-1,3-dicarboxamide (**7**)

1.3. <sup>1</sup>H NMR of *N, N'*-di-(2,3,4,6-tetra-*O*-acetyl-β-*D*-galactopyranosyl-1,2,3-triazol-4-ylmethyl amide)-*N''*-(2-azidoacetamido)-5-aminobenzene-1,3-dicarboxamide (**8**)

1.4. <sup>13</sup>C NMR of *N, N'*-di-(2,3,4,6-tetra-*O*-acetyl-β-*D*-galactopyranosyl-1,2,3-triazol-4-ylmethyl amide)-*N''*-(2-azidoacetamido)-5-aminobenzene-1,3-dicarboxamide (**8**)

1.5. <sup>1</sup>H NMR of 2-[2-(2-Propargyloxyethoxy)ethoxy]ethanol (**9**)

1.6. <sup>1</sup>H NMR of 2-(2-(2-Propargyloxyethoxy)ethoxy)ethyl-4-methylbenzenesulfonate (**10**)

1.7. <sup>1</sup>H NMR of *N, N'*-di-(2,3,4,6-tetra-*O*-acetyl-β-*D*-galactopyranosyl-1,2,3-triazol-4-ylmethyl amide)-*N''*-(2-4-((2-(2-(2-(4-methylbenzenesulfonate)ethoxy)ethoxy)ethoxy) methyl)-1*H*-1,2,3-triazol-1-yl)acetamido)-5-aminobenzene-1,3-dicarboxamide (**11**)

1.8. <sup>13</sup>C NMR of *N, N'*-di-(2,3,4,6-tetra-*O*-acetyl-β-*D*-galactopyranosyl-1,2,3-triazol-4-ylmethyl amide)-*N''*-(2-4-((2-(2-(2-(4-methylbenzenesulfonate)ethoxy)ethoxy)ethoxy) methyl)-1*H*-1,2,3-triazol-1-yl)acetamido)-5-aminobenzene-1,3-dicarboxamide (**11**)

1.9. <sup>1</sup>H NMR of *N, N'*-di-(2,3,4,6-tetra-*O*-acetyl-β-*D*-galactopyranosyl-1,2,3-triazol-4-ylmethylamide)-*N''*-(2-4-((2-(2-(2-azidoethoxy)ethoxy)ethoxy)methyl)-1*H*-1,2,3-triazol-1-yl)acetamido)-5-aminobenzene-1,3-dicarboxamide (**12**)

1.10. <sup>13</sup>C NMR of *N, N'*-di-(2,3,4,6-tetra-*O*-acetyl-β-*D*-galactopyranosyl-1,2,3-triazol-4-ylmethylamide)-*N''*-(2-4-((2-(2-(2-azidoethoxy)ethoxy)ethoxy)methyl)-1*H*-1,2,3-triazol-1-yl)acetamido)-5-aminobenzene-1,3-dicarboxamide (**12**)

1.11. <sup>1</sup>H NMR of *tert*-butyl (4,8,12,16-tetra-*aza*)(5,9,13,17-tetra-*oxo*)(4,8,12,16-tetra-*N*-propargyl)octadecanoate (**2**)

1.12.  $^{13}\text{C}$  NMR of tert-butyl (4,8,12,16-tetra-aza)(5,9,13,17-tetra-oxo)(4,8,12,16-tetra-N-propargyl) octadecanoate (**2**)

1.13.  $^1\text{H}$  NMR of Acetylated Tetravalent  $\beta$ -Peptoid Glycocluster (**13**)

1.14.  $^{13}\text{C}$  NMR of Acetylated Tetravalent  $\beta$ -Peptoid Glycocluster (**13**)

1.15.  $^1\text{H}$  NMR of Tetravalent  $\beta$ -Peptoid Glycocluster (**3**)

1.16.  $^{13}\text{C}$  NMR of Tetravalent  $\beta$ -Peptoid Glycocluster (**3**)

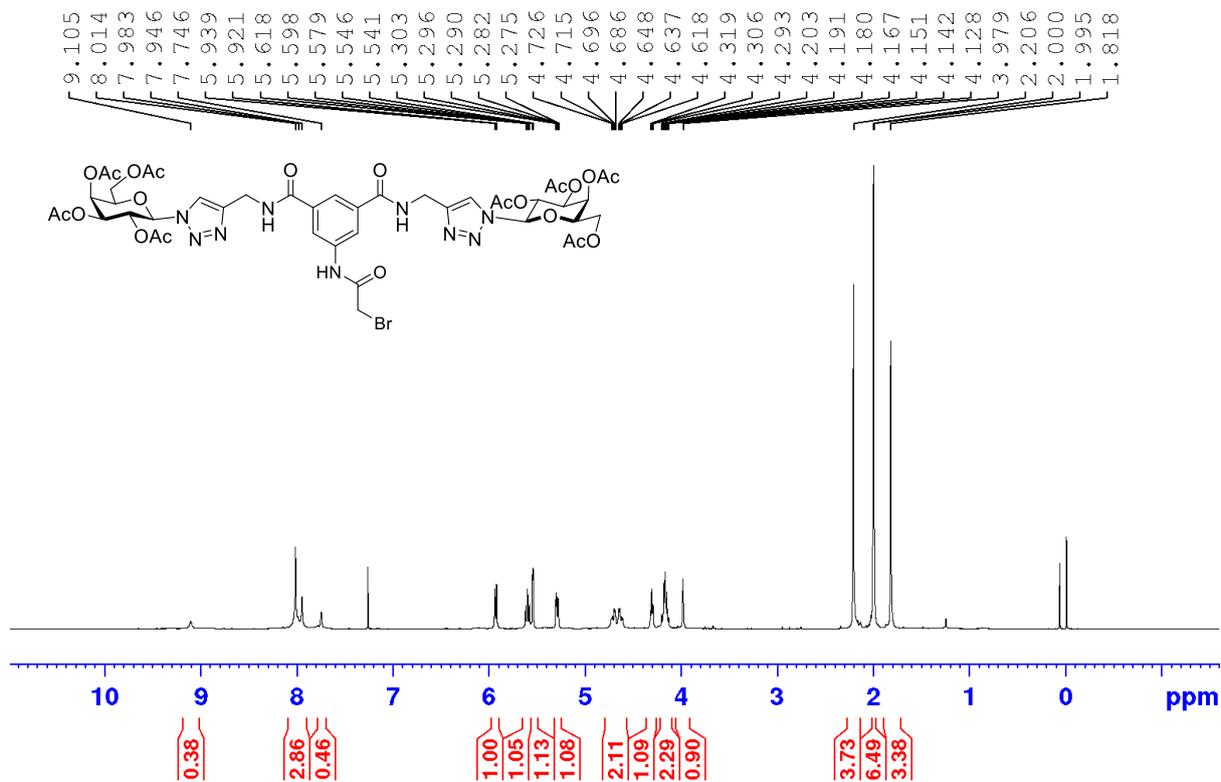
1.17.  $^1\text{H}$  NMR of 2,6,10,14-tetraoxo-3,7,11,15-tetrakis((1-(2,3,4,6-tetra-O-acetyl- $\beta$ -D-galactopyranosyl -1H-1,2,3-triazol-4-yl)methyl)-3,7,11,15-tetraazaoctadecan-18-oic acid (**14**)

1.18.  $^{13}\text{C}$  NMR of 2,6,10,14-tetraoxo-3,7,11,15-tetrakis((1-(2,3,4,6-tetra-O-acetyl- $\beta$ -D-galactopyranosyl -1H-1,2,3-triazol-4-yl)methyl)-3,7,11,15-tetraazaoctadecan-18-oic acid (**14**)

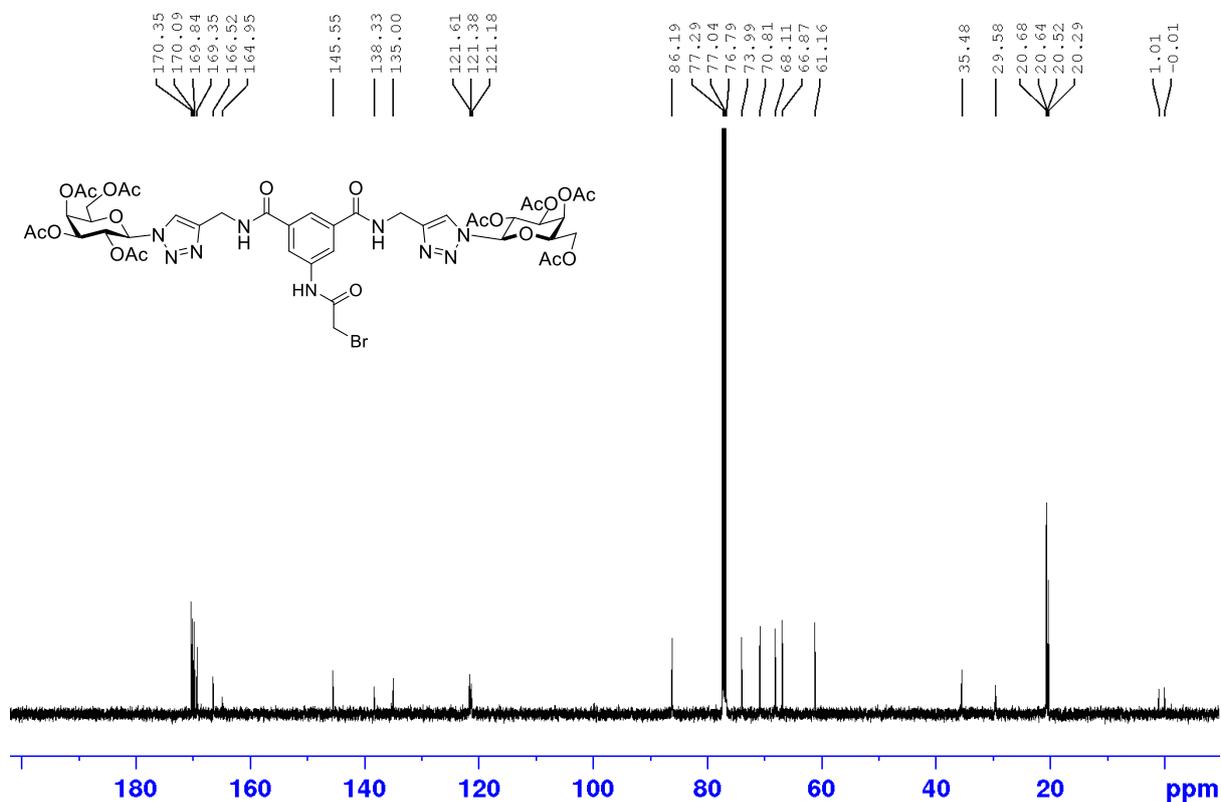
1.19.  $^1\text{H}$  NMR of 2,6,10,14-tetraoxo-3,7,11,15-tetrakis((1-( $\beta$ -D-galactopyranosyl -1H-1,2,3-triazol-4-yl)methyl)-3,7,11,15-tetraazaoctadecan-18-oic acid (**15**)

1.20.  $^{13}\text{C}$  NMR of 2,6,10,14-tetraoxo-3,7,11,15-tetrakis((1-( $\beta$ -D-galactopyranosyl -1H-1,2,3-triazol-4-yl)methyl)-3,7,11,15-tetraazaoctadecan-18-oic acid (**15**)

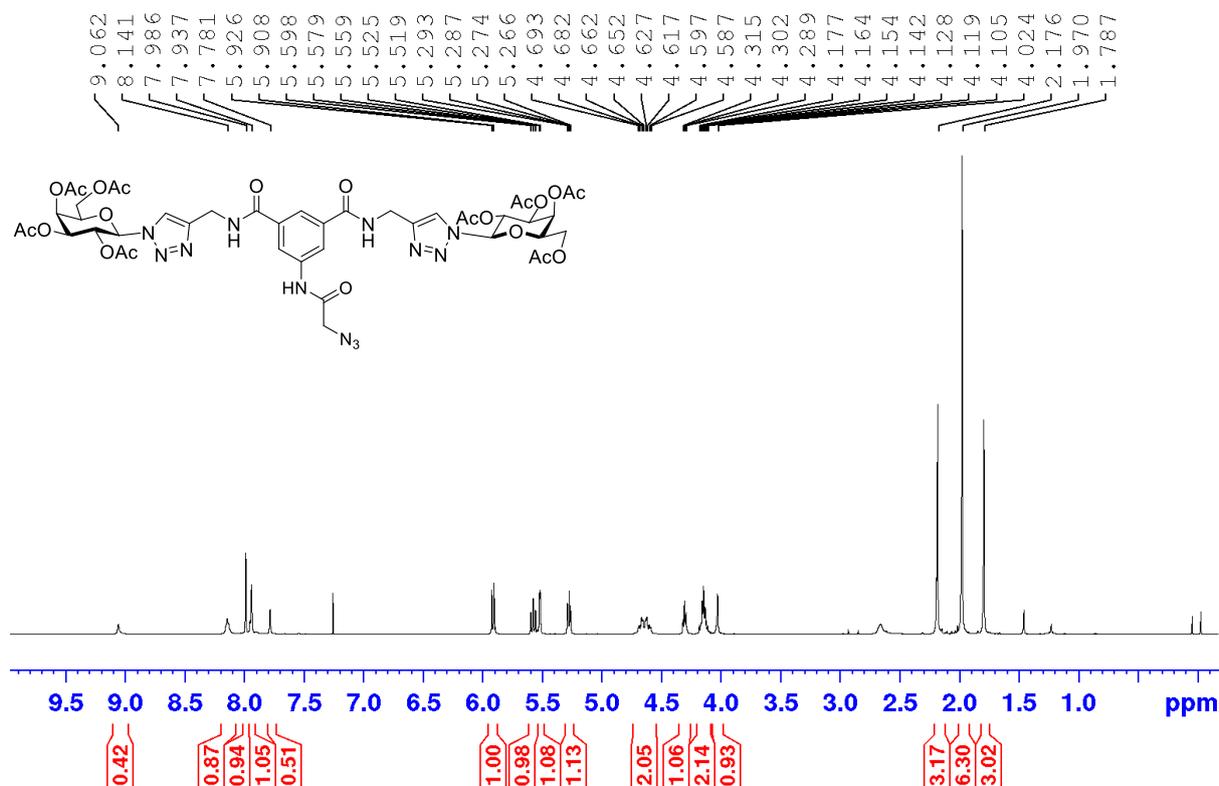
1.1.  $^1\text{H}$  NMR of *N, N'*-di-(2,3,4,6-tetra-*O*-acetyl- $\beta$ -*D*-galactopyranosyl-1,2,3-triazol-4-ylmethylamide)-*N''*-(2-bromoacetamido)-5-aminobenzene-1,3-dicarboxamide (**7**)



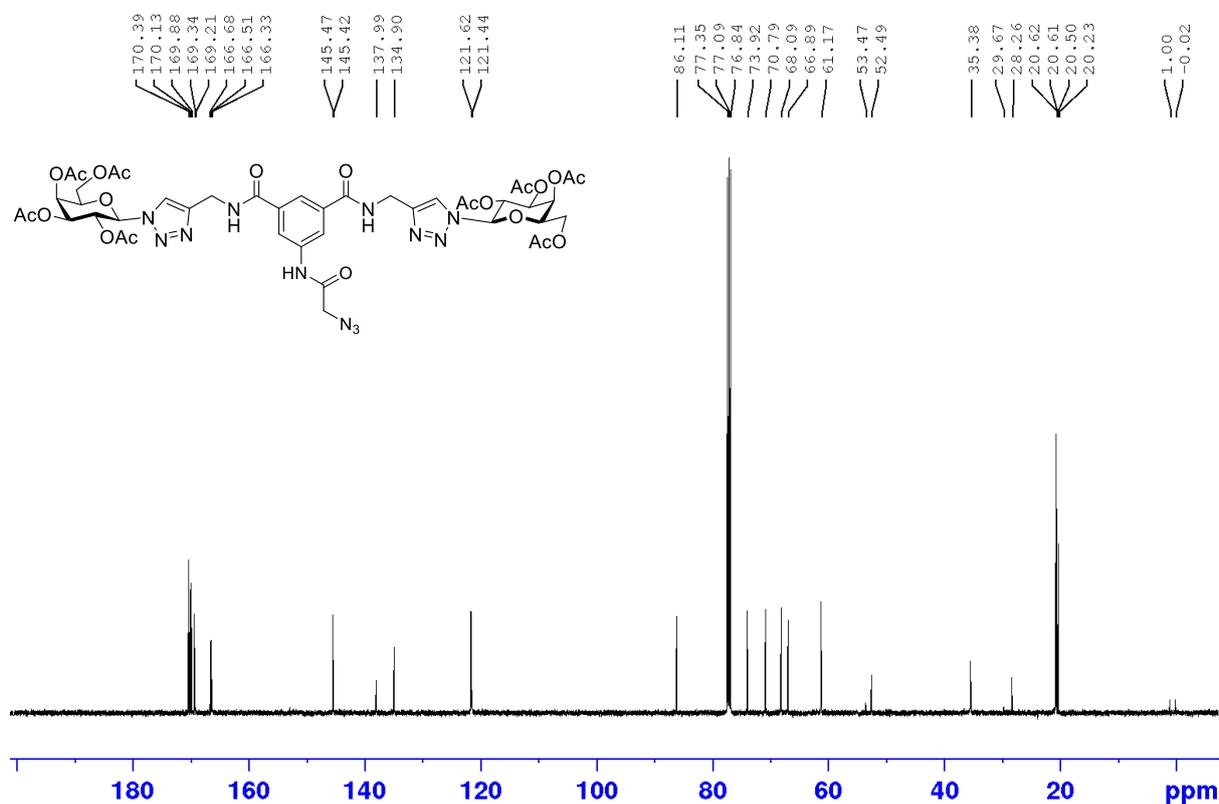
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1.3.  $^1\text{H}$  NMR of *N,N'*-di-(2,3,4,6-tetra-*O*-acetyl- $\beta$ -D-galactopyranosyl-1,2,3-triazol-4-ylmethyl amide)-*N''*-(2-azidoacetamido)-5-aminobenzene-1,3-dicarboxamide (**8**)

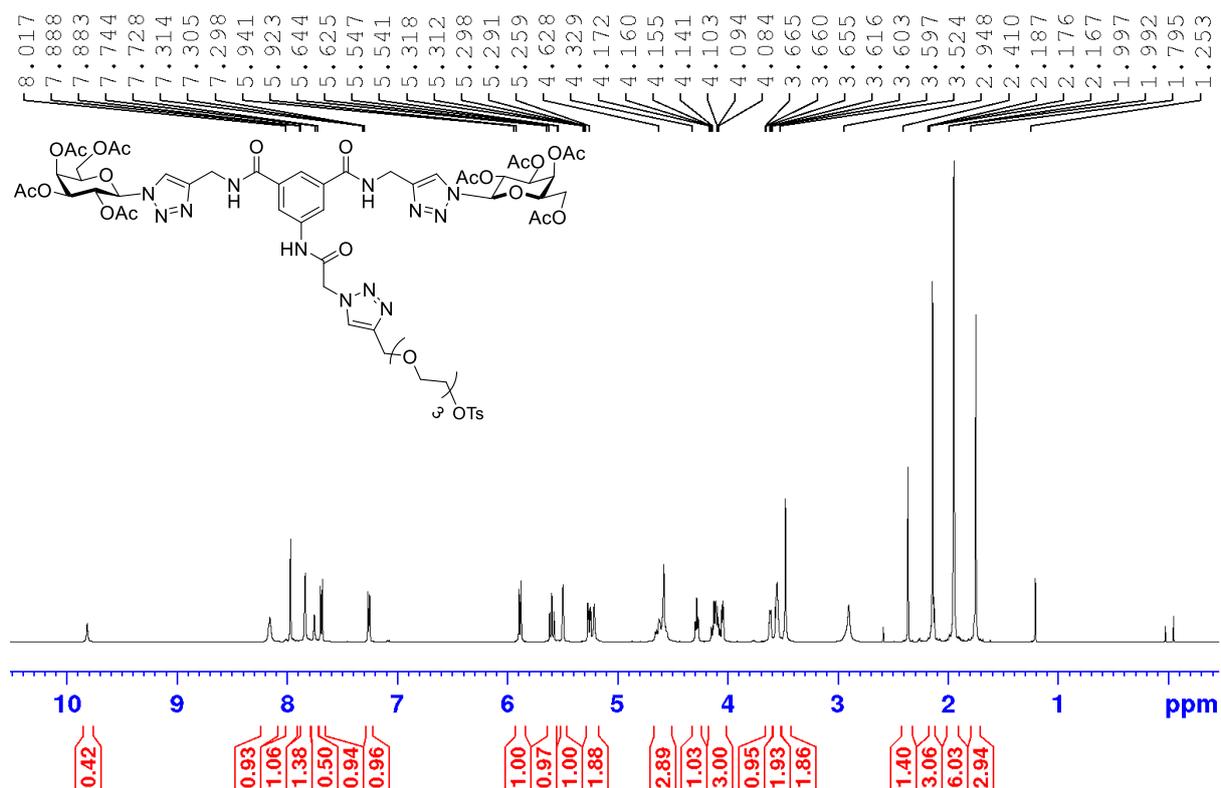


1.4.  $^{13}\text{C}$  NMR of *N,N'*-di-(2,3,4,6-tetra-*O*-acetyl- $\beta$ -D-galactopyranosyl-1,2,3-triazol-4-ylmethyl amide)-*N''*-(2-azidoacetamido)-5-aminobenzene-1,3-dicarboxamide (**8**)

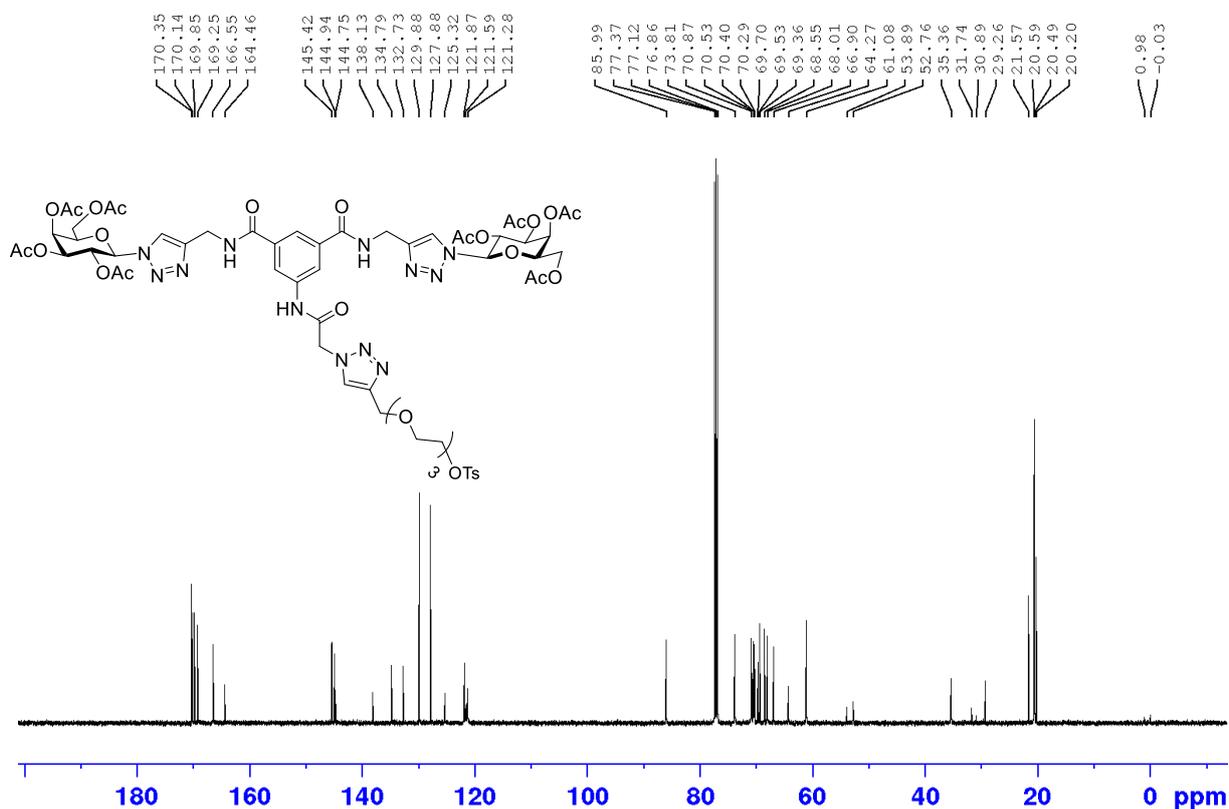




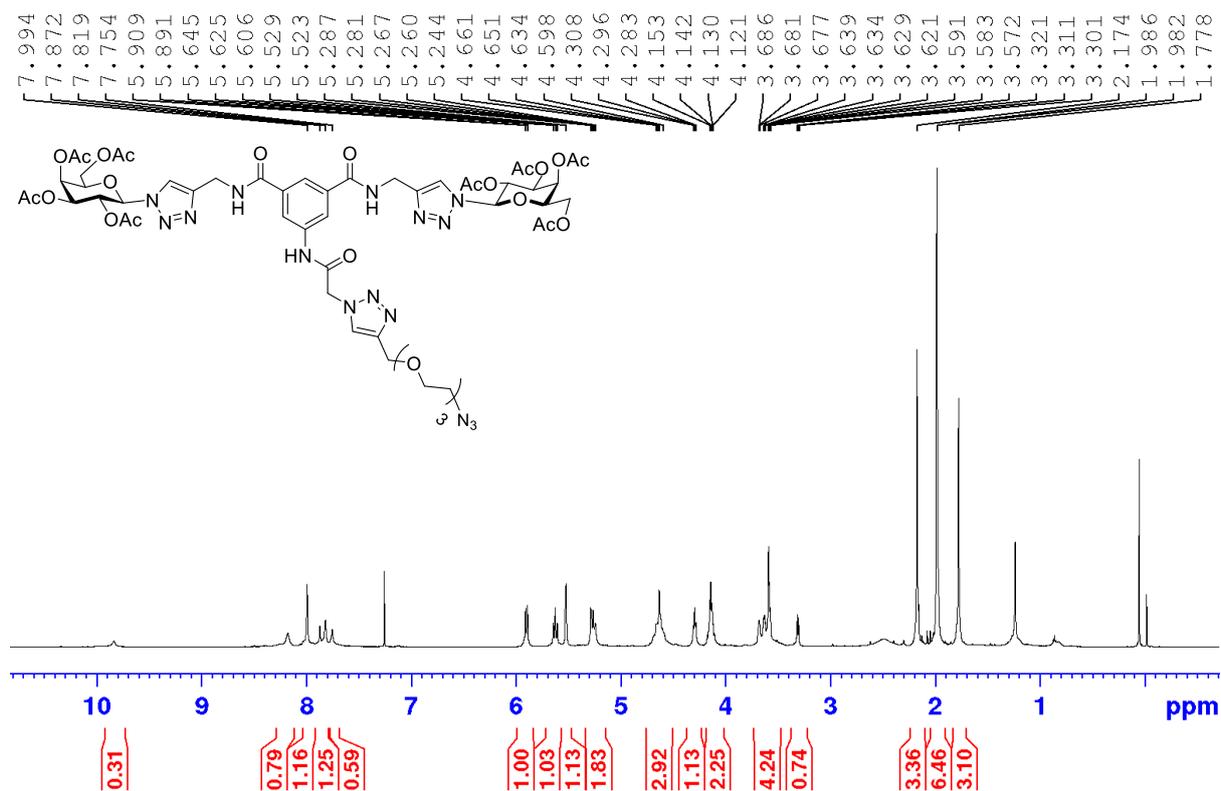
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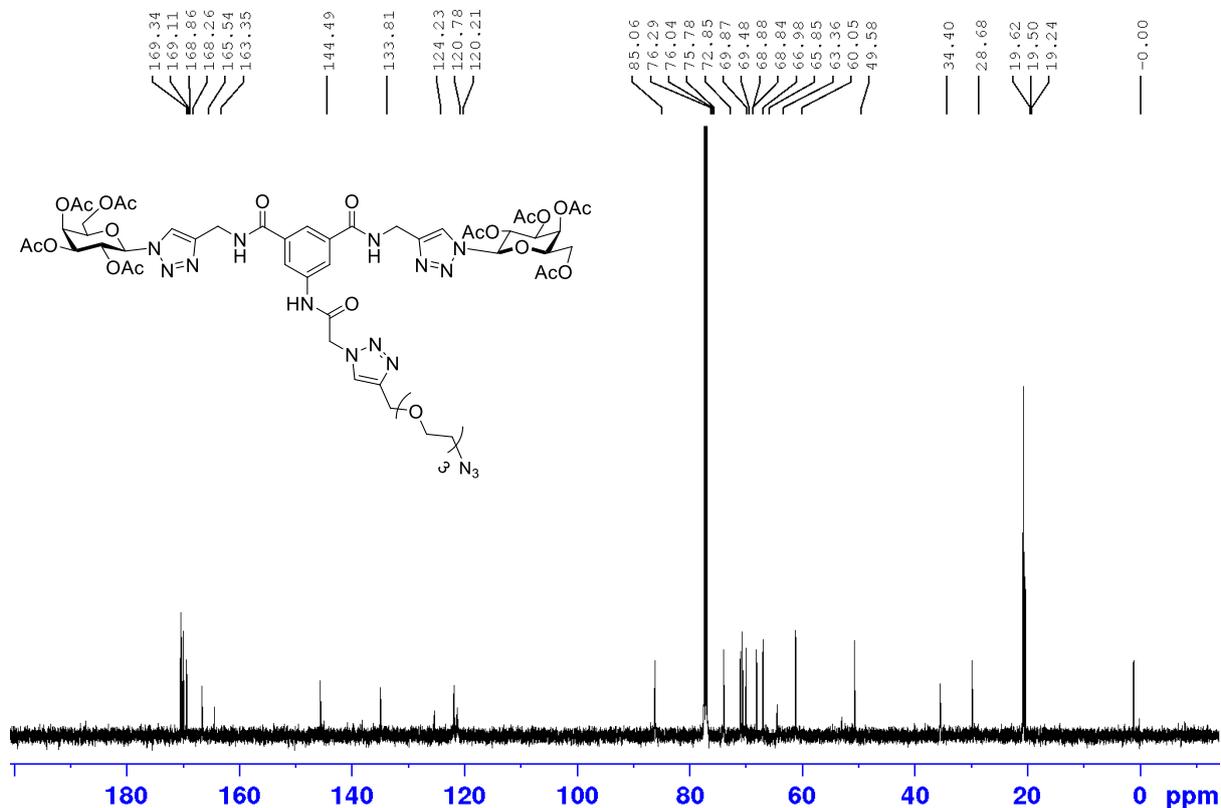
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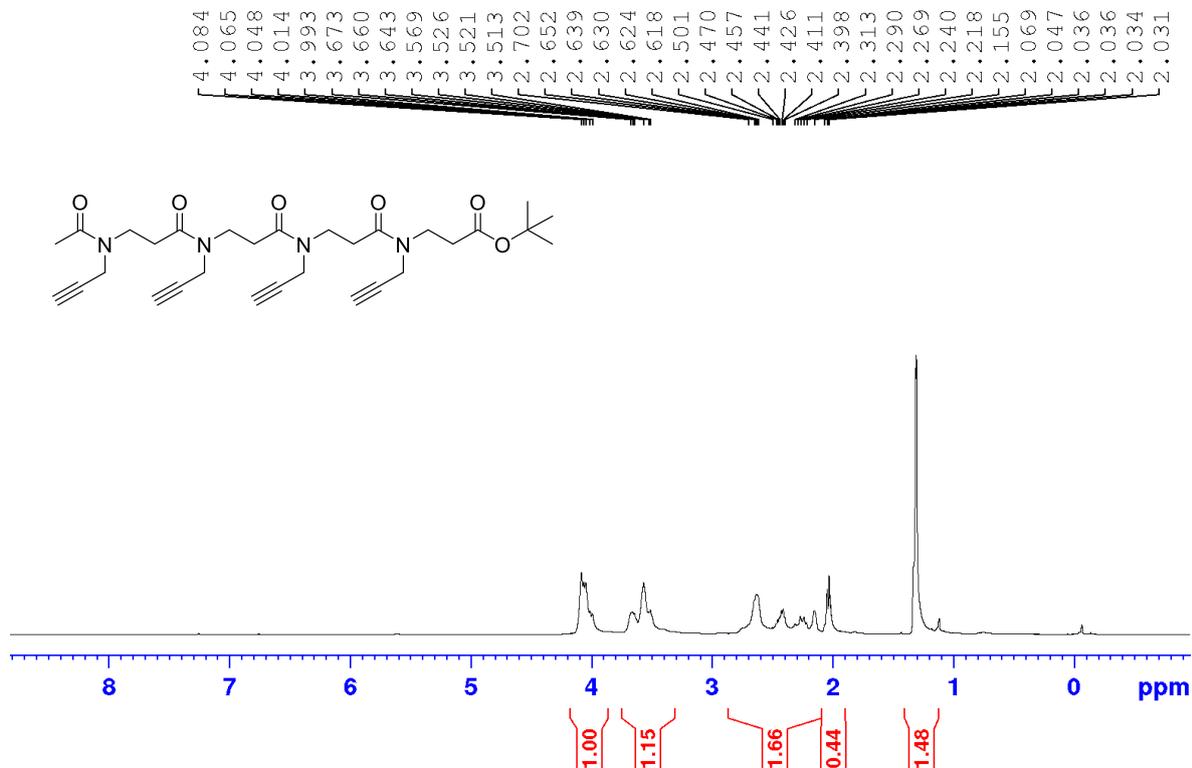
1.9.  $^1\text{H}$  NMR of *N, N'*-di-(2,3,4,6-tetra-*O*-acetyl- $\beta$ -*D*-galactopyranosyl-1,2,3-triazol-4-yl)methylamide)-*N''*-(2-4-((2-(2-(2-azidoethoxy)ethoxy)ethoxy)methyl)-1*H*-1,2,3-triazol-1-yl)acetamido)-5-aminobenzene-1,3-dicarboxamide (**12**)



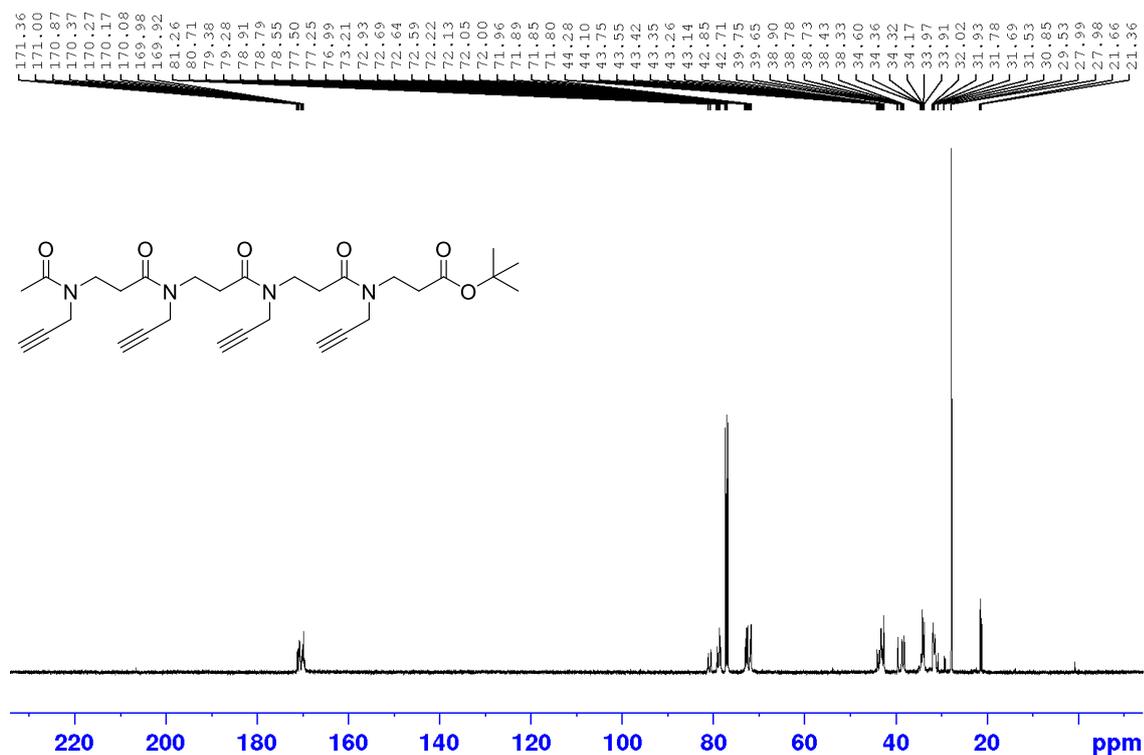
1.10.  $^{13}\text{C}$  NMR of *N, N'*-di-(2,3,4,6-tetra-*O*-acetyl- $\beta$ -*D*-galactopyranosyl-1,2,3-triazol-4-yl)methylamide)-*N''*-(2-4-((2-(2-(2-azidoethoxy)ethoxy)ethoxy)methyl)-1*H*-1,2,3-triazol-1-yl)acetamido)-5-aminobenzene-1,3-dicarboxamide (**12**)



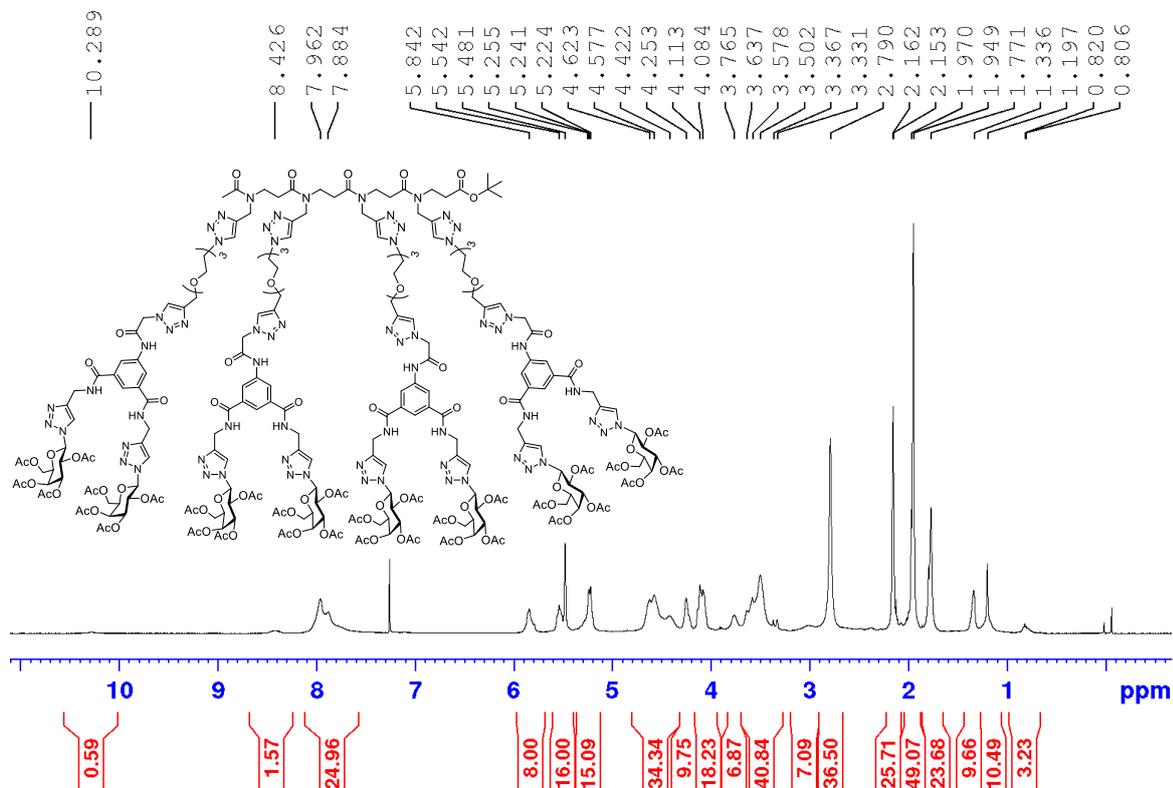
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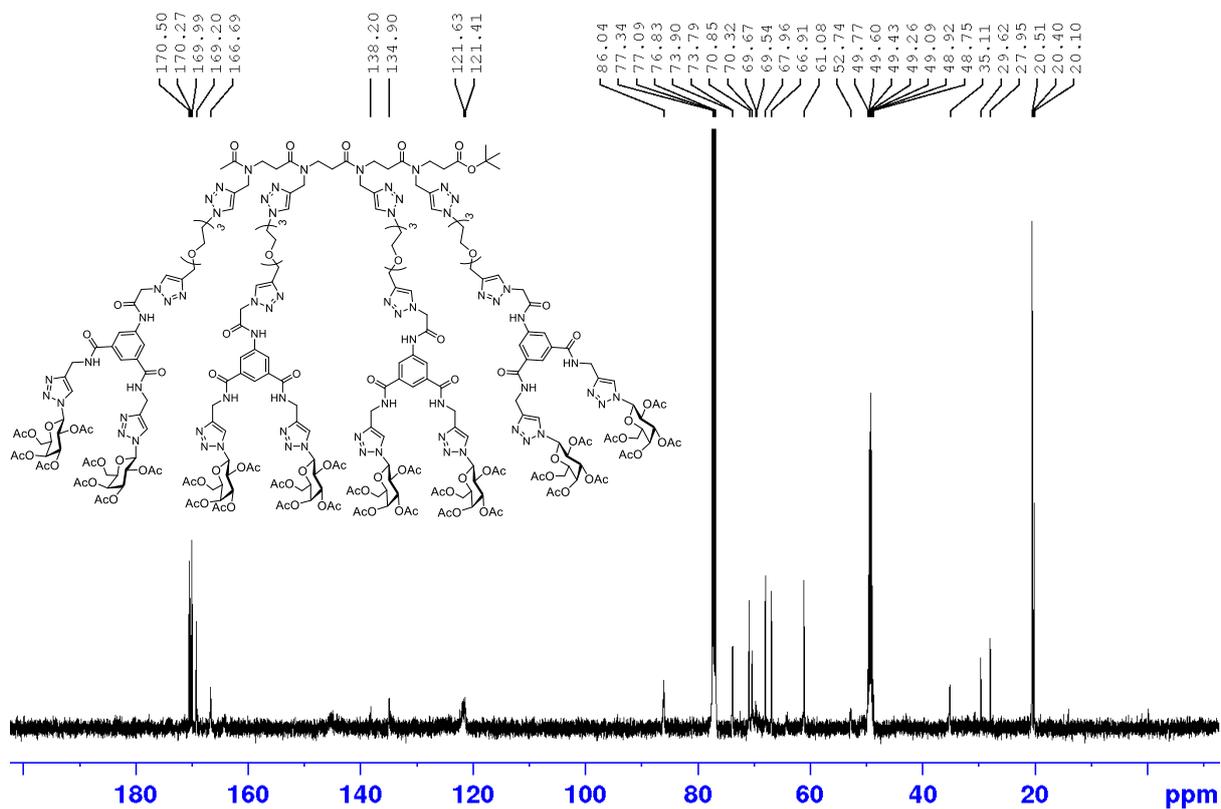
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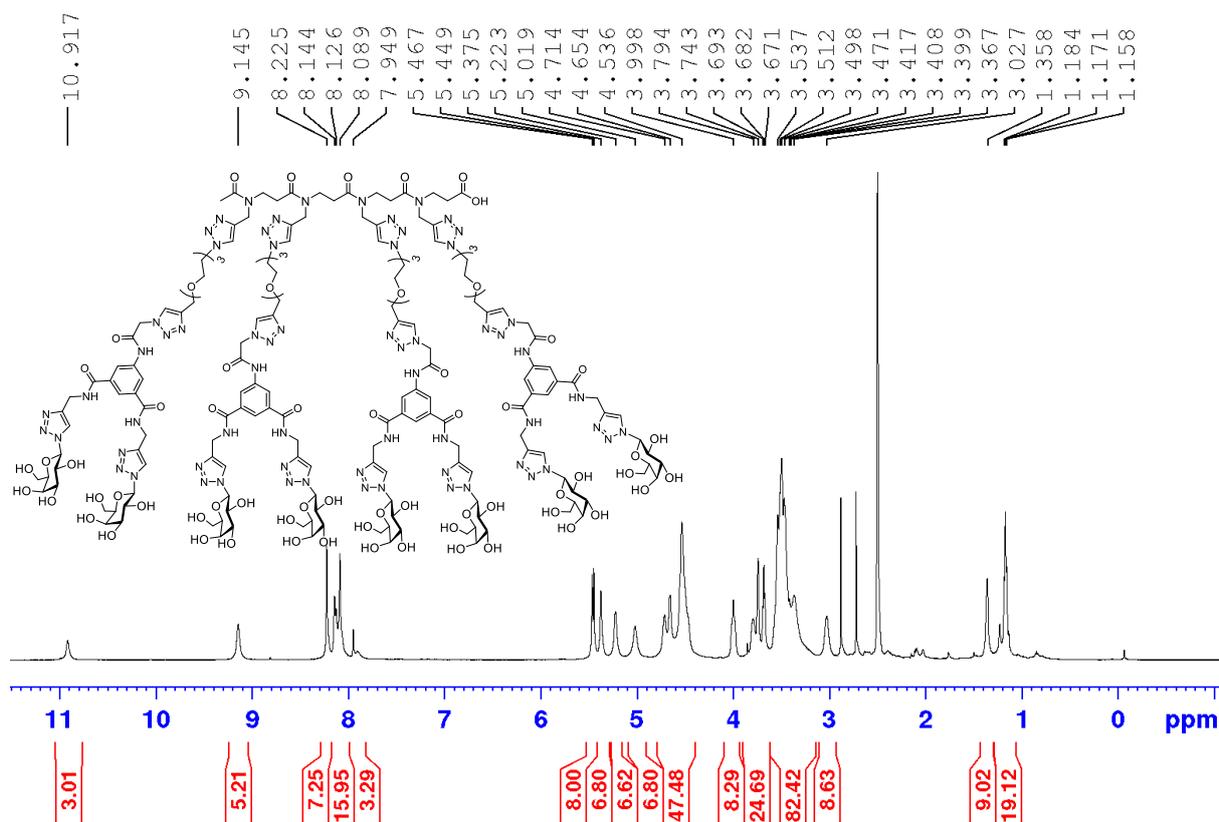
1.13. <sup>1</sup>H NMR of Acetylated Tetravalent β-Peptoid Glycocluster (13)



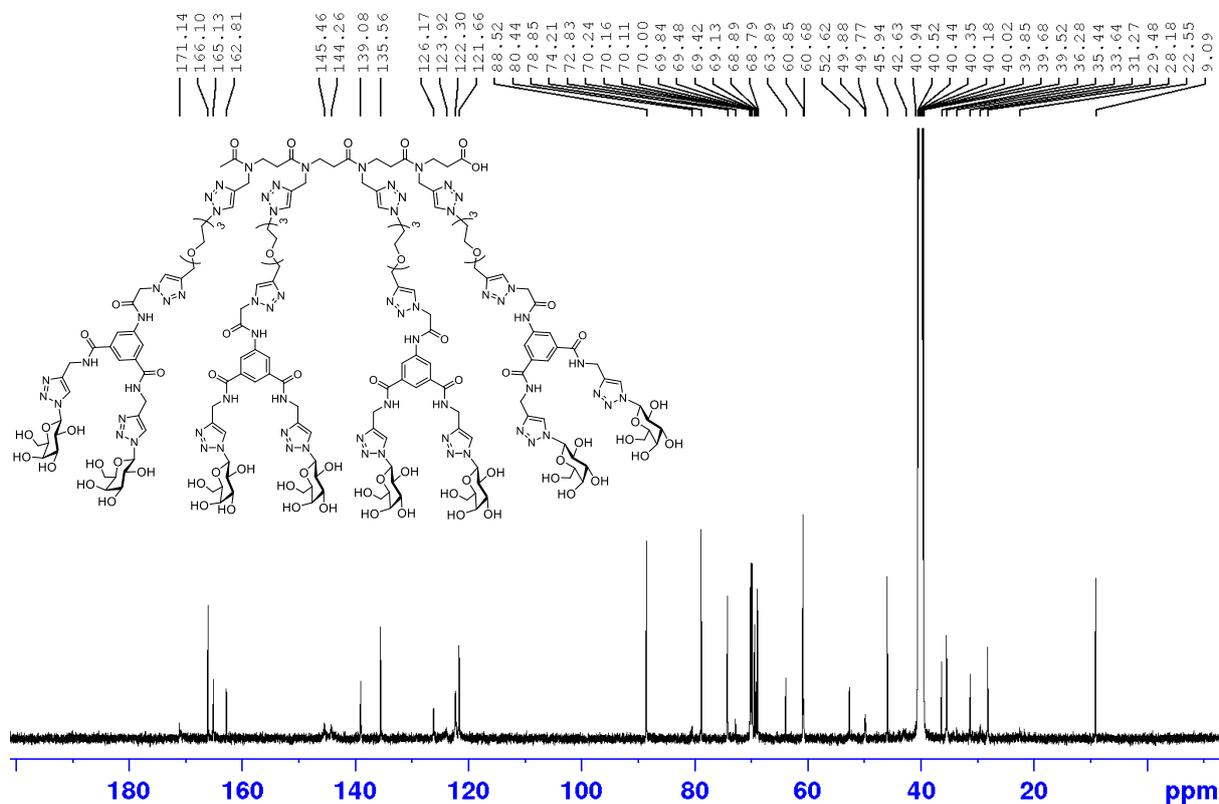
1.14. <sup>13</sup>C NMR of Acetylated Tetravalent β-Peptoid Glycocluster (13)



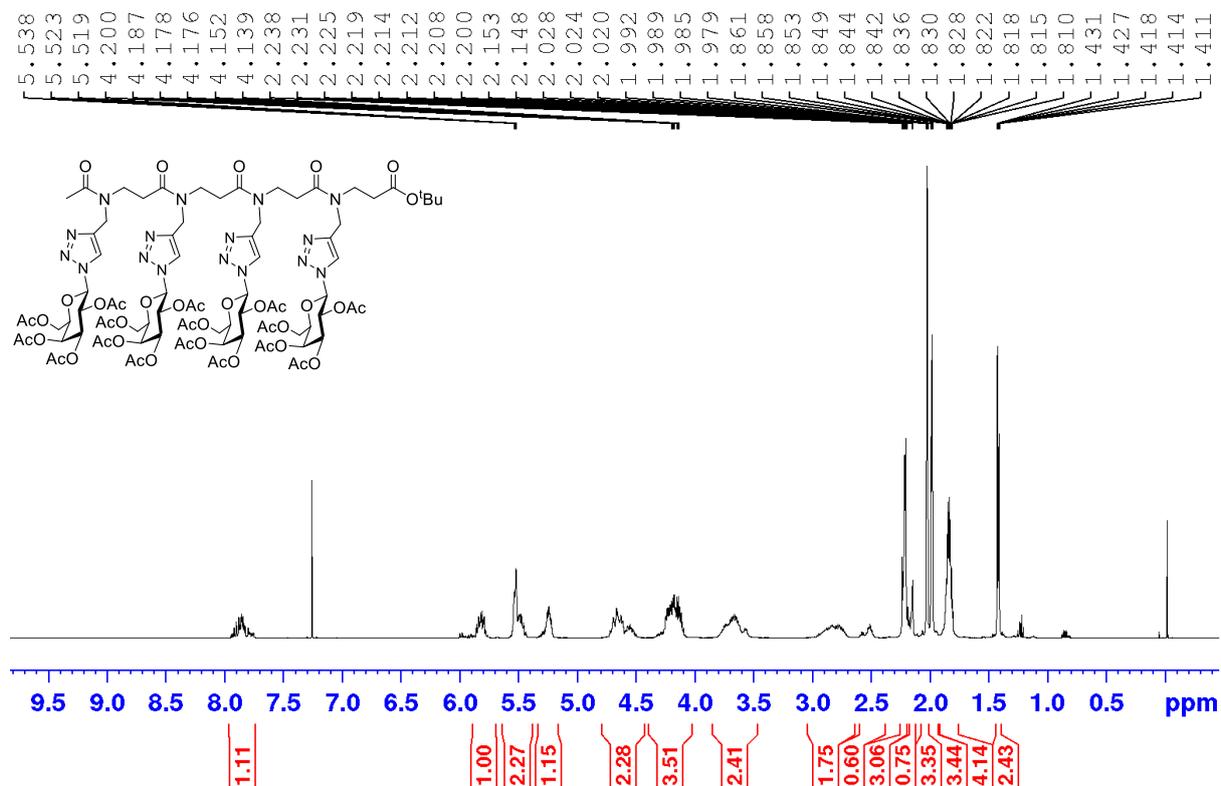
### 1.15. <sup>1</sup>H NMR of Tetravalent β-Peptoid Glycocluster (3)



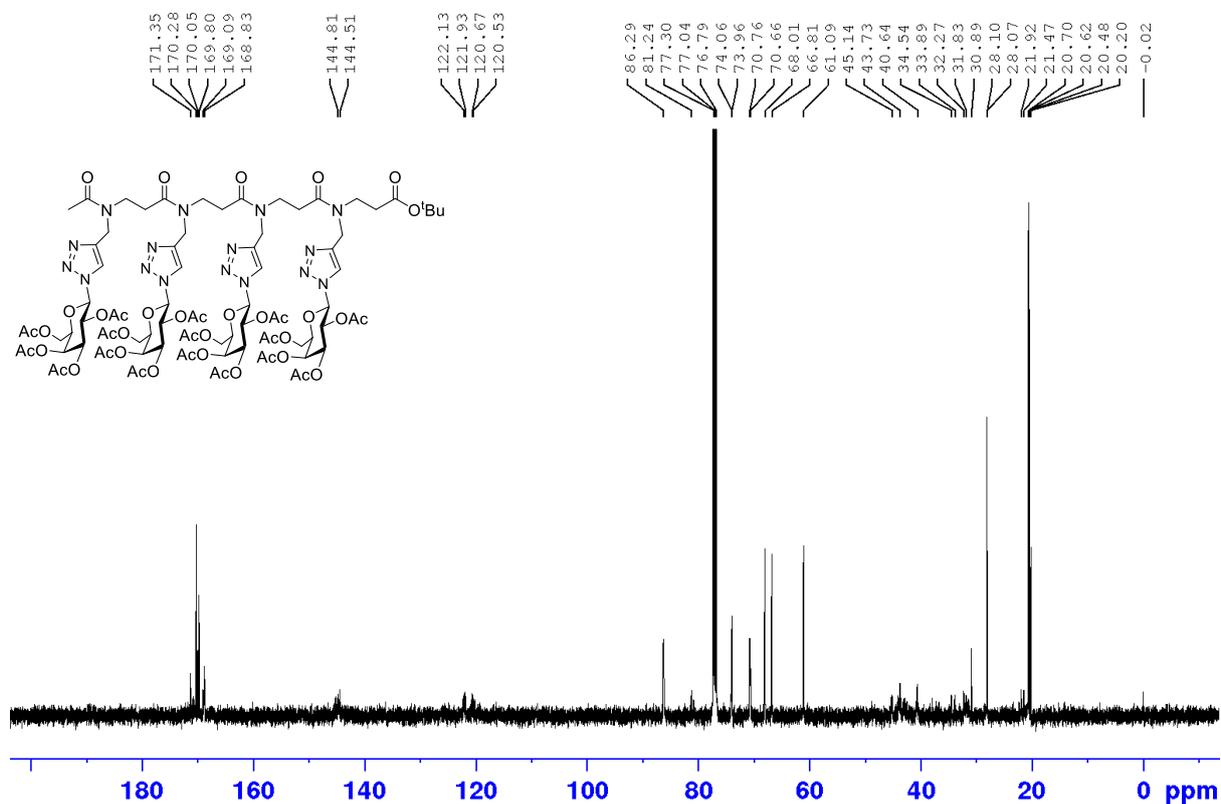
### 1.16. <sup>13</sup>C NMR of Tetravalent β-Peptoid Glycocluster (3)



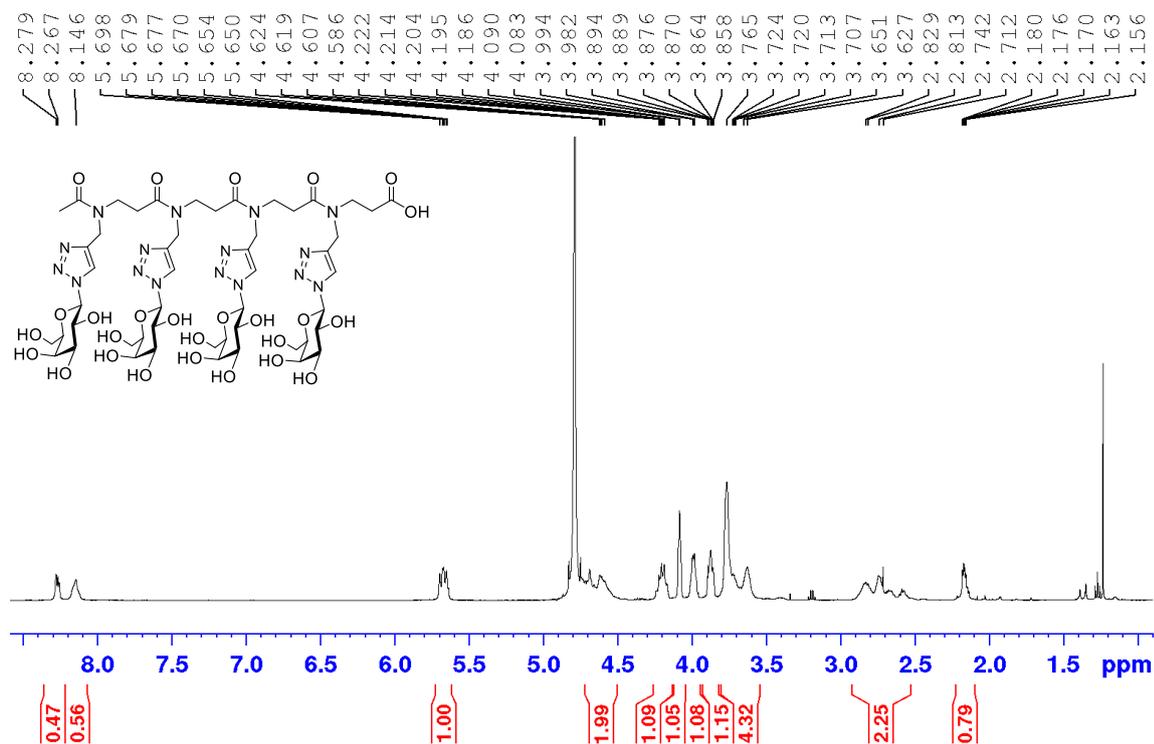
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