



# wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 9, 2024 – 10:27 am GMT

PDB ID : 5IEE  
Title : Murine endoplasmic reticulum alpha-glucosidase II with 1-deoxynojirimycin  
Authors : Caputo, A.T.; Roversi, P.; Alonzi, D.S.; Kiappes, J.L.; Zitzmann, N.  
Deposited on : 2016-02-25  
Resolution : 1.92 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 3.0  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.003 (Gargrove)  
Density-Fitness : 1.0.11  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.39

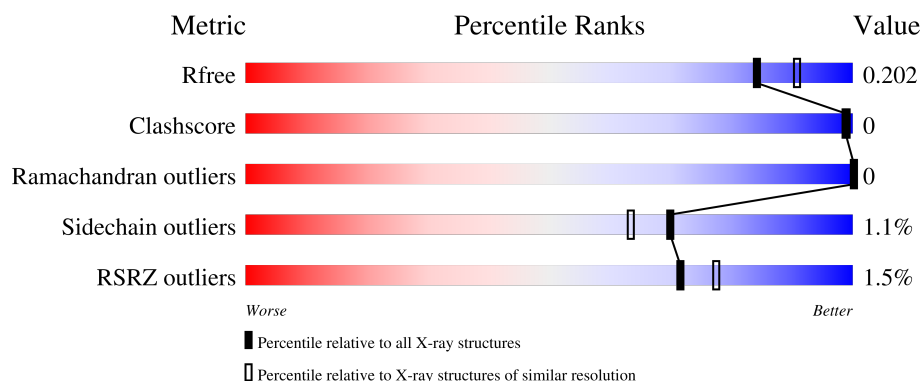
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.92 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	164625	1028 (1.92-1.92)
Clashscore	180529	1100 (1.92-1.92)
Ramachandran outliers	177936	1087 (1.92-1.92)
Sidechain outliers	177891	1087 (1.92-1.92)
RSRZ outliers	164620	1028 (1.92-1.92)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	934	<div> <div style="width: 100%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">%</span> </div> <div style="width: 90%; height: 10px; background-color: green; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">90%</span> </div> <div style="width: 8%; height: 10px; background-color: grey; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">8%</span> </div> </div>
2	B	88	<div> <div style="width: 3%; height: 10px; background-color: red; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">3%</span> </div> <div style="width: 100%; height: 10px; background-color: green; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">100%</span> </div> </div>
3	C	2	<div> <div style="width: 50%; height: 10px; background-color: green; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">50%</span> </div> <div style="width: 50%; height: 10px; background-color: yellow; position: relative;"> <span style="position: absolute; top: -10px; left: 0;">50%</span> </div> </div>

## 2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 16328 atoms, of which 7536 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Neutral alpha-glucosidase AB.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	857	Total	C	H	N	O	S	0	13	0
			13728	4477	6745	1207	1269	30			

- Molecule 2 is a protein called Glucosidase 2 subunit beta.

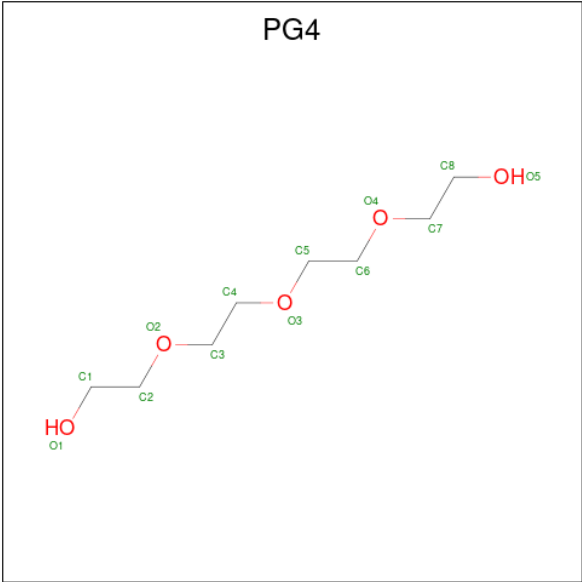
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	88	Total	C	H	N	O	S	0	0	0
			1231	399	568	106	148	10			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



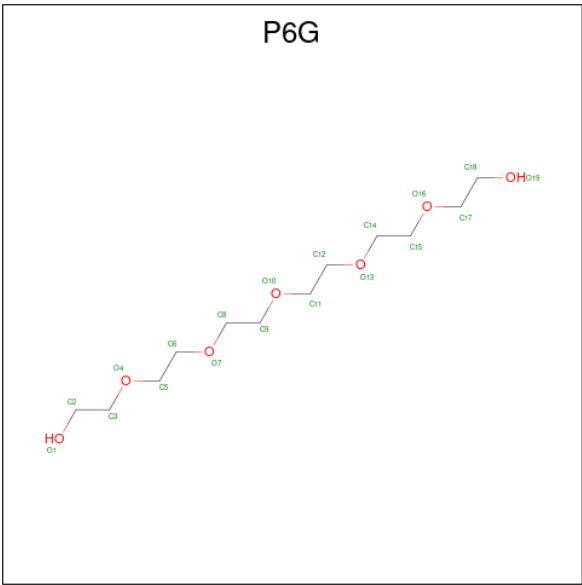
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	2	Total	C	H	N	O		0	0	0
			53	16	25	2	10				

- Molecule 4 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C<sub>8</sub>H<sub>18</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			17	5	9	3		
4	A	1	Total	C	H	O	0	0
			20	6	11	3		
4	A	1	Total	C	H	O	0	0
			31	8	18	5		

- Molecule 5 is HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula: C<sub>12</sub>H<sub>26</sub>O<sub>7</sub>).



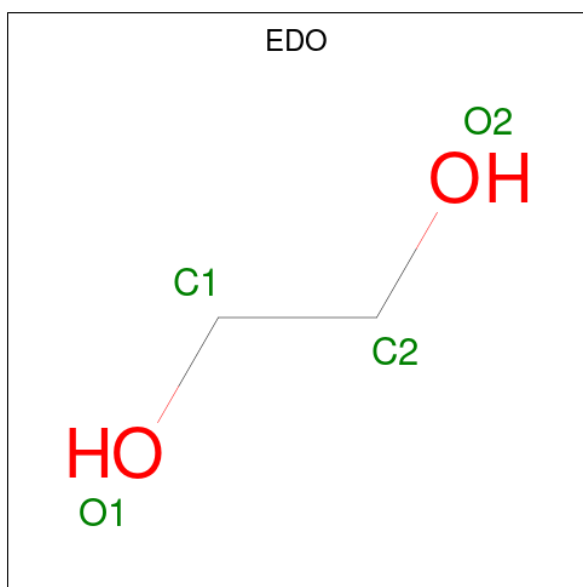
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			45	12	26	7		

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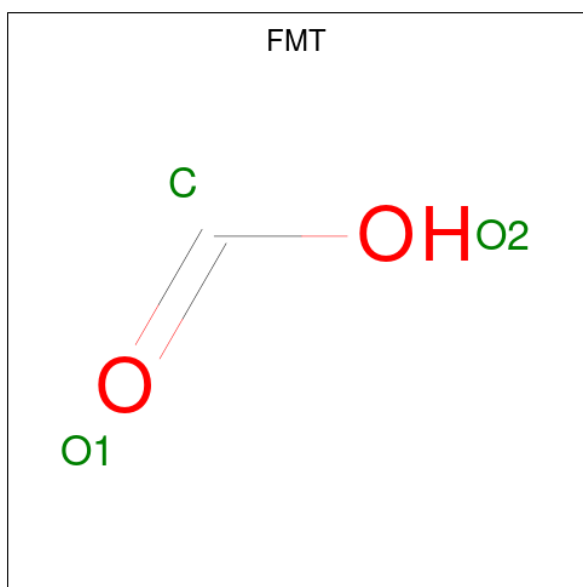
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			45	12	26	7		
5	A	1	Total	C	H	O	0	0
			30	9	16	5		
5	A	1	Total	C	H	O	0	0
			45	12	26	7		

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



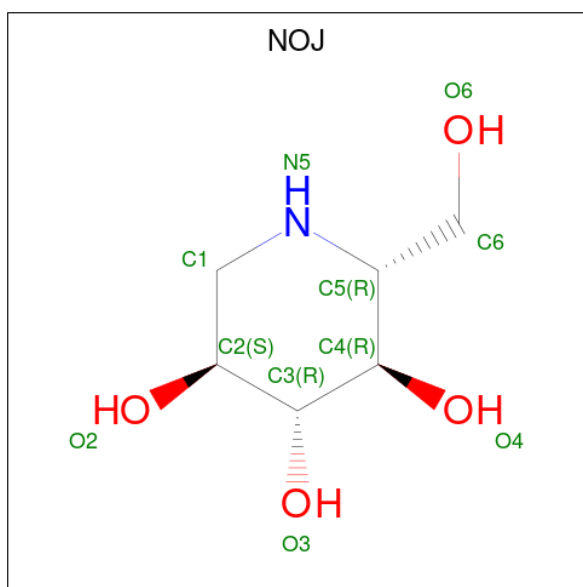
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	B	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 7 is FORMIC ACID (three-letter code: FMT) (formula:  $CH_2O_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	H	O	0	0
			4	1	1	2		
7	A	1	Total	C	H	O	0	0
			4	1	1	2		
7	A	1	Total	C	H	O	0	0
			5	1	2	2		
7	A	1	Total	C	H	O	0	0
			5	1	2	2		
7	A	1	Total	C	H	O	0	0
			5	1	2	2		
7	A	1	Total	C	H	O	0	0
			4	1	1	2		
7	A	1	Total	C	H	O	0	0
			5	1	2	2		
7	A	1	Total	C	H	O	0	0
			5	1	2	2		
7	B	1	Total	C	H	O	0	0
			5	1	2	2		

- Molecule 8 is 1-DEOXYNOJIRIMYCIN (three-letter code: NOJ) (formula: C<sub>6</sub>H<sub>13</sub>NO<sub>4</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	H	N	O	0	0
			24	6	13	1	4		

- Molecule 9 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	B	2	Total	Ca	0	0
			2	2		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	848	Total	O	0	0
			848	848		
10	B	102	Total	O	0	0
			102	102		





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 <sub>1</sub> 2 <sub>1</sub> 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	103.48Å 173.71Å 62.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	88.90 – 1.92 88.90 – 1.92	Depositor EDS
% Data completeness (in resolution range)	99.9 (88.90-1.92) 99.9 (88.90-1.92)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.24	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.56 (at 1.92Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, $R_{free}$	0.164 , 0.196 0.164 , 0.202	Depositor DCC
$R_{free}$ test set	4314 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.2	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.44 , 64.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	16328	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NOJ, P6G, PG4, NAG, CA, FMT, EDO

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.50	0/7242	0.68	0/9856
2	B	0.49	0/677	0.64	0/920
All	All	0.50	0/7919	0.68	0/10776

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6983	6745	6746	6	0
2	B	663	568	568	0	0
3	C	28	25	25	0	0
4	A	30	38	38	0	0
5	A	71	94	94	0	0
6	A	20	30	30	1	0
6	B	4	6	6	0	0
7	A	27	15	15	0	0
7	B	3	2	1	0	0
8	A	11	13	13	0	0
9	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	A	848	0	0	1	0
10	B	102	0	0	0	0
All	All	8792	7536	7536	6	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 0.

The worst 5 of 6 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:927:LEU:HD23	10:A:1225:HOH:O	1.96	0.66
1:A:248:TRP:O	1:A:259:LYS:HE3	1.95	0.65
1:A:432:LEU:HD22	1:A:477[A]:HIS:CD2	2.42	0.53
1:A:320:SER:O	1:A:627:PHE:HA	2.10	0.51
1:A:802:GLN:HG3	6:A:1011:EDO:C2	2.42	0.50

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	864/934 (92%)	838 (97%)	26 (3%)	0	100	100
2	B	86/88 (98%)	85 (99%)	1 (1%)	0	100	100
All	All	950/1022 (93%)	923 (97%)	27 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	761/810 (94%)	752 (99%)	9 (1%)	67	59
2	B	77/77 (100%)	77 (100%)	0	100	100
All	All	838/887 (94%)	829 (99%)	9 (1%)	70	63

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	685	TRP
1	A	706	ARG
1	A	424	ASN
1	A	446	ASP
1	A	637	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

2 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	NAG	C	1	3,1	14,14,15	0.33	0	17,19,21	0.89	1 (5%)
3	NAG	C	2	3	14,14,15	0.28	0	17,19,21	0.59	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	C	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	C	2	3	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	NAG	C1-O5-C5	2.05	114.96	112.19

There are no chirality outliers.

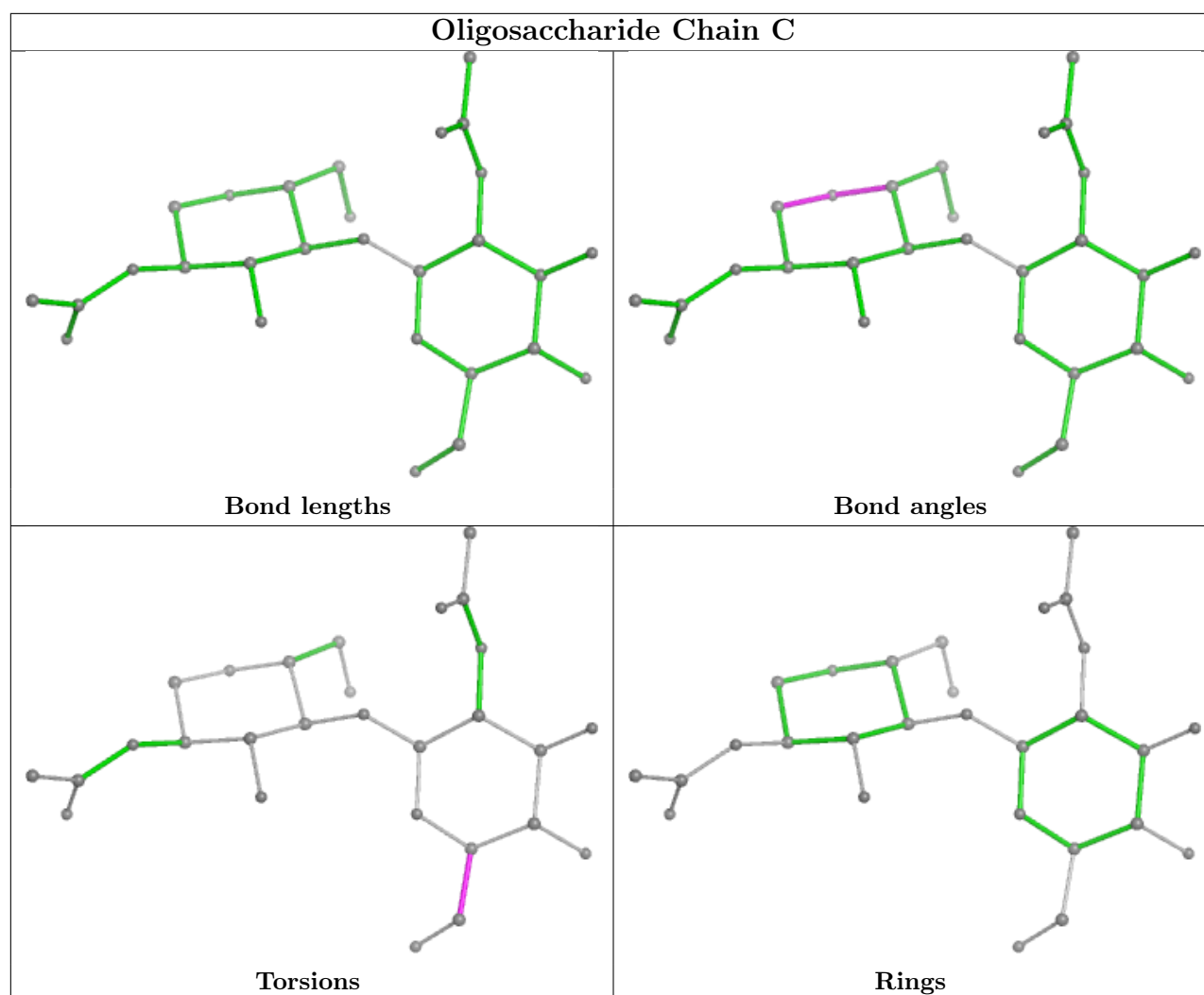
All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	2	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



## 5.6 Ligand geometry [i](#)

Of 26 ligands modelled in this entry, 2 are monoatomic - leaving 24 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
4	PG4	A	1005	-	12,12,12	0.25	0	11,11,11	0.33	0
7	FMT	A	1019	-	2,2,2	2.08	1 (50%)	1,1,1	1.23	0
6	EDO	A	1013	-	3,3,3	0.56	0	2,2,2	0.24	0
8	NOJ	A	1023	-	11,11,11	0.26	0	13,15,15	1.10	1 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
7	FMT	A	1014	-	2,2,2	0.91	0	1,1,1	1.15	0
7	FMT	A	1021	-	2,2,2	1.24	0	1,1,1	1.06	0
7	FMT	A	1016	-	2,2,2	1.56	1 (50%)	1,1,1	1.00	0
6	EDO	B	203	-	3,3,3	0.47	0	2,2,2	0.38	0
4	PG4	A	1003	-	7,7,12	0.19	0	6,6,11	0.18	0
7	FMT	A	1018	-	2,2,2	1.16	0	1,1,1	1.11	0
5	P6G	A	1008	-	13,13,18	0.13	0	12,12,17	0.19	0
7	FMT	A	1017	-	2,2,2	1.36	0	1,1,1	1.04	0
7	FMT	A	1022	-	2,2,2	1.72	1 (50%)	1,1,1	1.11	0
6	EDO	A	1011	-	3,3,3	0.34	0	2,2,2	0.83	0
5	P6G	A	1024	-	18,18,18	0.27	0	17,17,17	0.30	0
5	P6G	A	1006	-	18,18,18	0.24	0	17,17,17	0.30	0
6	EDO	A	1012	-	3,3,3	0.60	0	2,2,2	0.26	0
6	EDO	A	1009	-	3,3,3	0.60	0	2,2,2	0.33	0
6	EDO	A	1010	-	3,3,3	0.68	0	2,2,2	0.29	0
7	FMT	B	204	-	2,2,2	1.13	0	1,1,1	1.05	0
5	P6G	A	1007	-	18,18,18	0.11	0	17,17,17	0.20	0
4	PG4	A	1004	-	8,8,12	0.25	0	7,7,11	0.30	0
7	FMT	A	1020	-	2,2,2	1.40	0	1,1,1	0.96	0
7	FMT	A	1015	-	2,2,2	0.91	0	1,1,1	1.43	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PG4	A	1005	-	-	6/10/10/10	-
6	EDO	B	203	-	-	1/1/1/1	-
4	PG4	A	1003	-	-	0/5/5/10	-
5	P6G	A	1008	-	-	2/11/11/16	-
5	P6G	A	1007	-	-	0/16/16/16	-
5	P6G	A	1006	-	-	7/16/16/16	-
4	PG4	A	1004	-	-	3/6/6/10	-
6	EDO	A	1012	-	-	1/1/1/1	-
6	EDO	A	1009	-	-	1/1/1/1	-
6	EDO	A	1013	-	-	1/1/1/1	-
6	EDO	A	1011	-	-	1/1/1/1	-
6	EDO	A	1010	-	-	1/1/1/1	-
5	P6G	A	1024	-	-	6/16/16/16	-
8	NOJ	A	1023	-	-	0/2/19/19	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	1019	FMT	O1-C	2.69	1.36	1.22
7	A	1022	FMT	O2-C	2.40	1.40	1.28
7	A	1016	FMT	O2-C	2.21	1.39	1.28

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	1023	NOJ	C1-N5-C5	3.46	117.10	109.61

There are no chirality outliers.

5 of 30 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1004	PG4	O2-C3-C4-O3
5	A	1024	P6G	O10-C11-C12-O13
5	A	1024	P6G	O13-C14-C15-O16
5	A	1006	P6G	O7-C8-C9-O10
5	A	1024	P6G	O16-C17-C18-O19

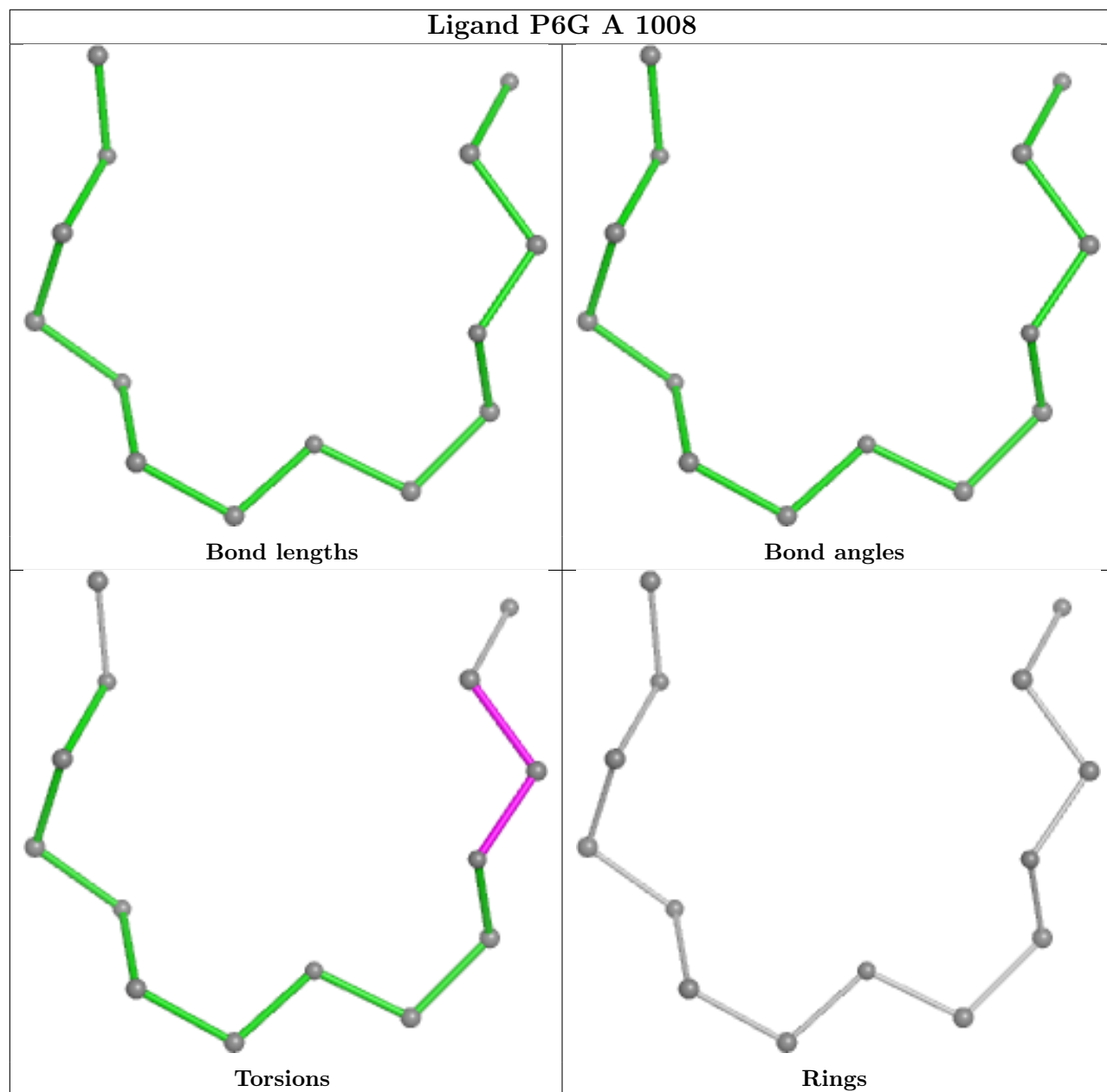
There are no ring outliers.

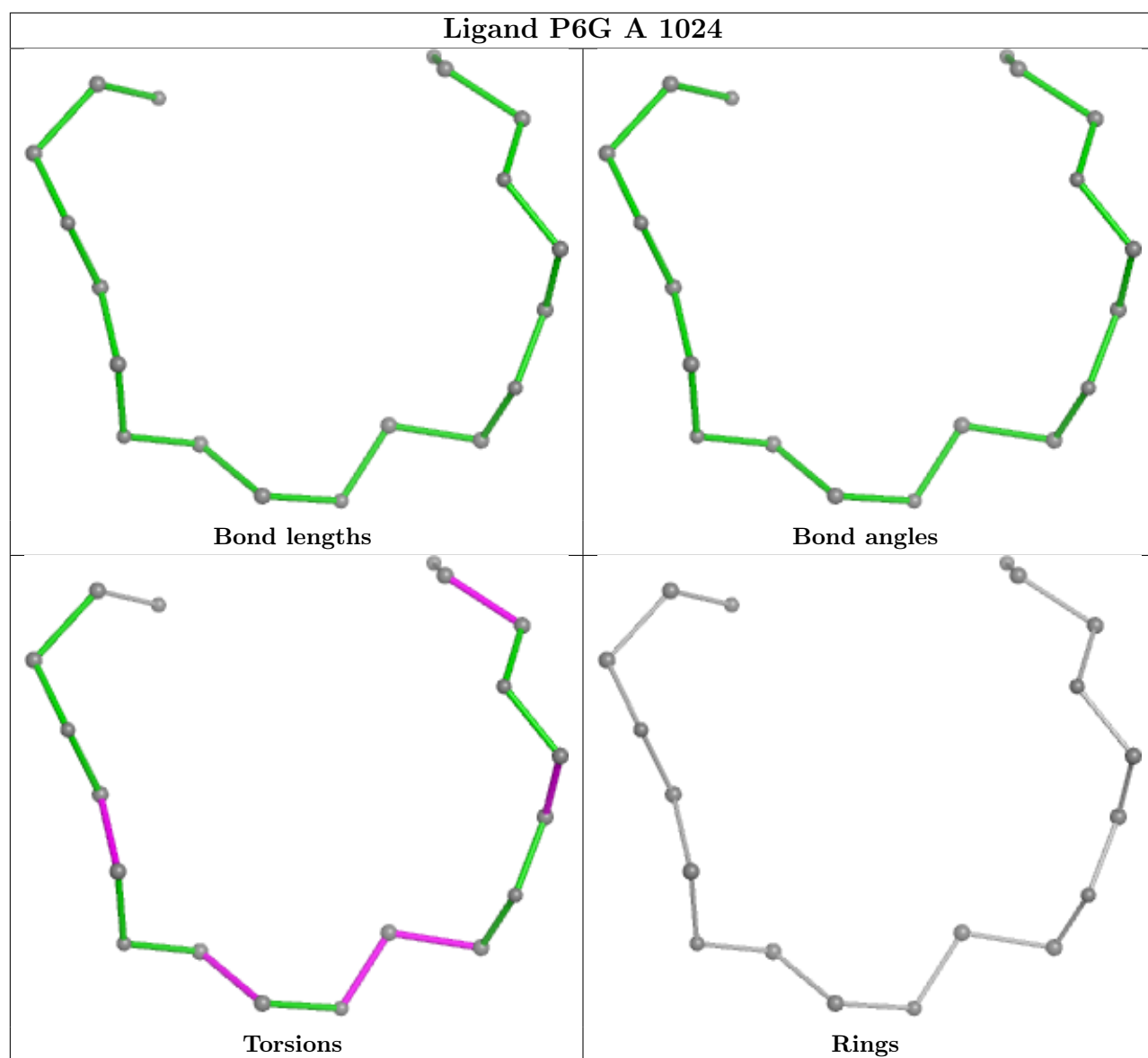
1 monomer is involved in 1 short contact:

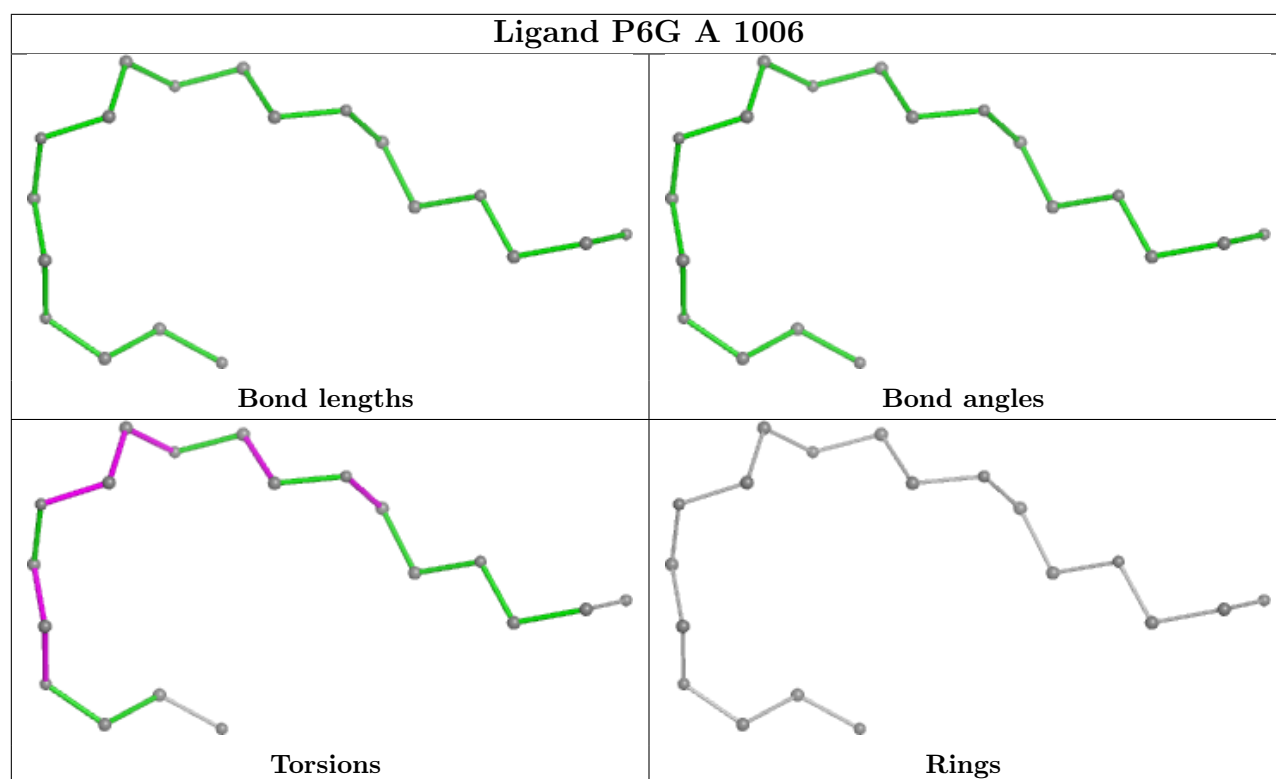
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1011	EDO	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	857/934 (91%)	-0.54	11 (1%) 74 80	11, 23, 44, 75	12 (1%)
2	B	88/88 (100%)	0.08	3 (3%) 48 55	20, 31, 53, 76	0
All	All	945/1022 (92%)	-0.49	14 (1%) 71 77	11, 23, 45, 76	12 (1%)

The worst 5 of 14 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	523[A]	TRP	3.7
1	A	350	ASN	3.7
2	B	30	PHE	3.3
1	A	54	ARG	3.1
2	B	48	PHE	3.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

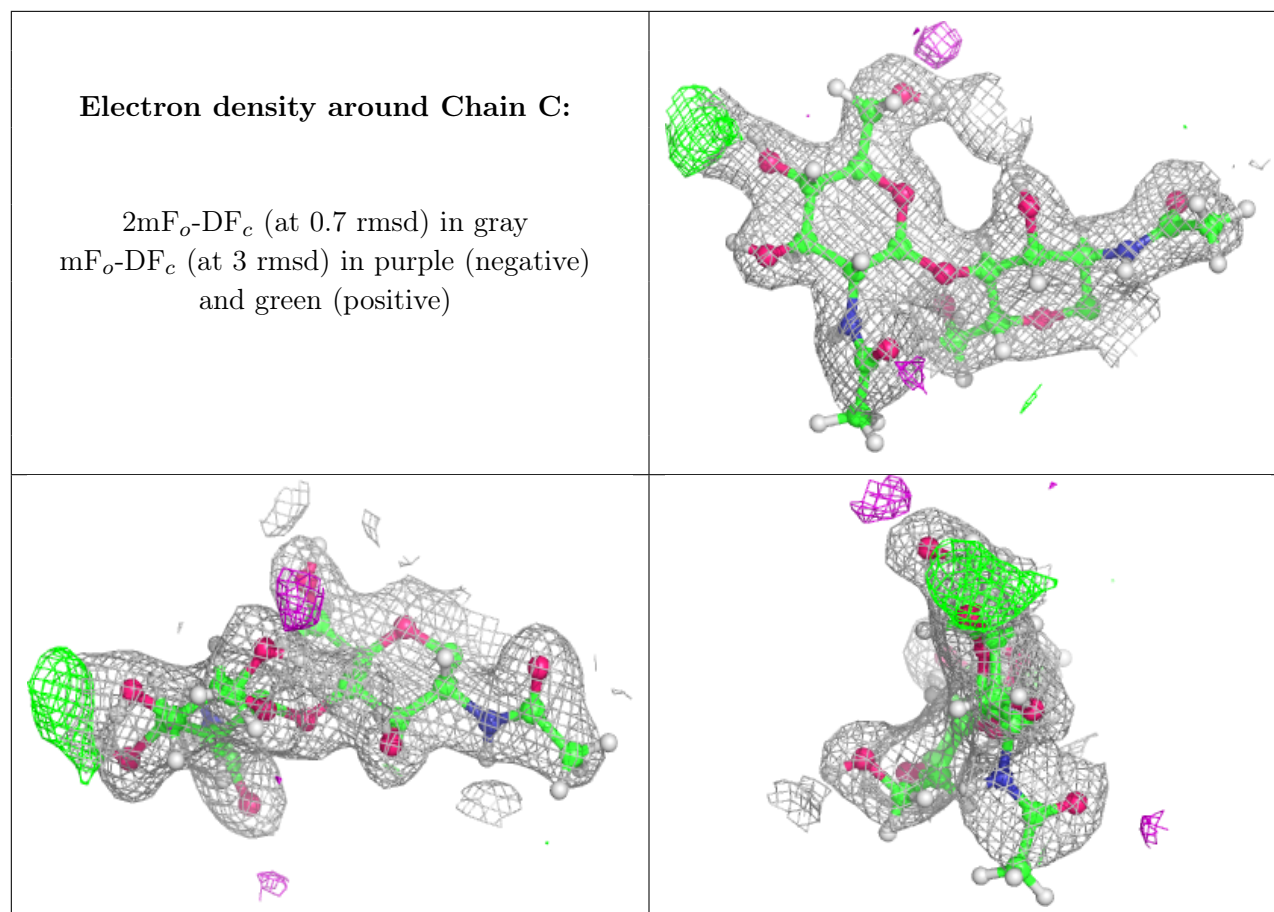
There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	NAG	C	2	14/15	0.92	0.08	34,38,43,45	0
3	NAG	C	1	14/15	0.97	0.05	24,27,33,33	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

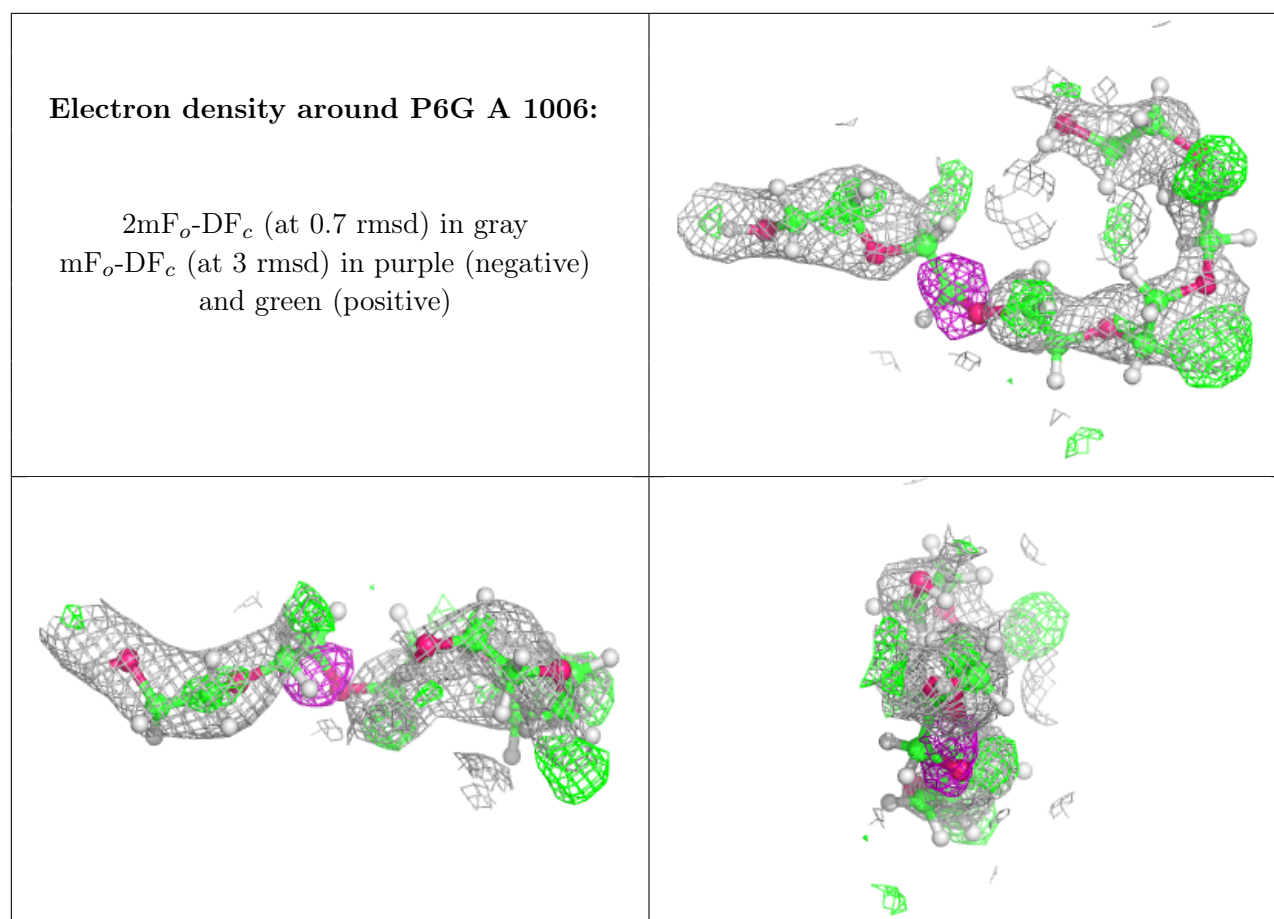
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
7	FMT	A	1014	3/3	0.69	0.30	68,70,70,71	0
7	FMT	A	1020	3/3	0.70	0.25	71,71,72,73	0
4	PG4	A	1004	9/13	0.73	0.21	47,60,63,64	0
5	P6G	A	1006	19/19	0.73	0.22	55,76,78,79	0
7	FMT	A	1022	3/3	0.79	0.16	30,30,35,38	0
6	EDO	A	1010	4/4	0.83	0.16	41,48,52,52	0
7	FMT	A	1016	3/3	0.84	0.26	53,53,54,54	0
7	FMT	A	1017	3/3	0.84	0.17	38,45,49,50	0
7	FMT	A	1021	3/3	0.85	0.17	64,64,66,66	0
5	P6G	A	1024	19/19	0.86	0.17	52,58,63,64	0
6	EDO	A	1009	4/4	0.86	0.14	48,52,60,60	0
4	PG4	A	1005	13/13	0.86	0.14	48,51,59,60	0

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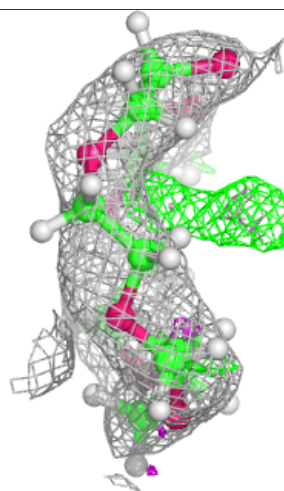
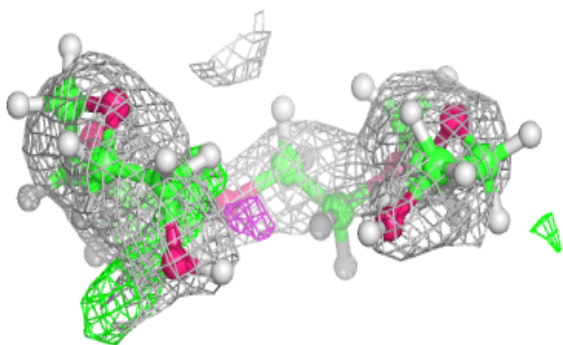
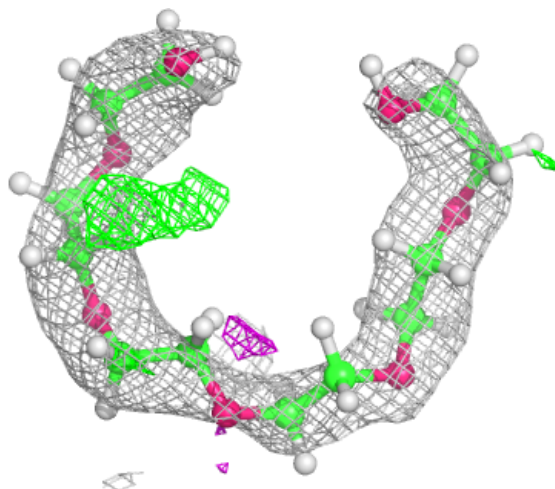
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
7	FMT	B	204	3/3	0.86	0.18	45,45,47,54	0
4	PG4	A	1003	8/13	0.87	0.14	41,41,50,52	0
5	P6G	A	1008	14/19	0.87	0.15	39,48,63,63	0
6	EDO	A	1013	4/4	0.89	0.12	53,55,55,55	0
7	FMT	A	1019	3/3	0.89	0.23	54,54,54,56	0
6	EDO	A	1011	4/4	0.90	0.11	49,52,59,60	0
6	EDO	B	203	4/4	0.90	0.14	53,55,57,57	0
7	FMT	A	1018	3/3	0.90	0.12	56,57,59,60	0
6	EDO	A	1012	4/4	0.90	0.14	46,47,51,51	0
7	FMT	A	1015	3/3	0.91	0.18	26,29,32,38	0
5	P6G	A	1007	19/19	0.92	0.10	36,44,53,56	0
8	NOJ	A	1023	11/11	0.98	0.04	17,18,21,21	0
9	CA	B	201	1/1	0.99	0.01	25,25,25,25	0
9	CA	B	202	1/1	1.00	0.02	22,22,22,22	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



**Electron density around P6G A 1024:**

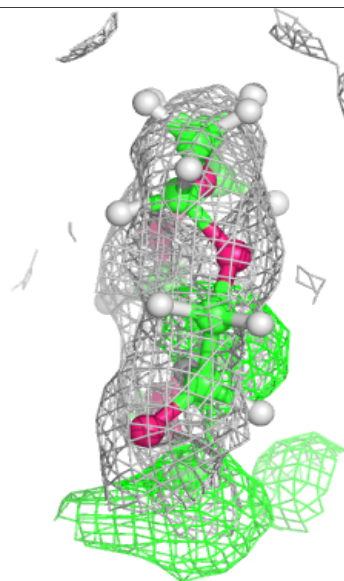
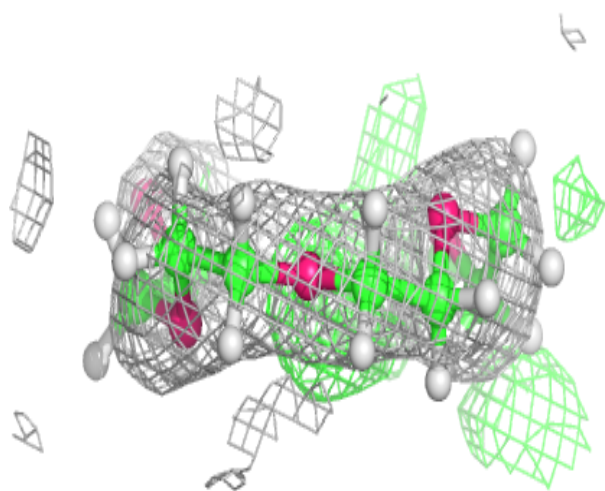
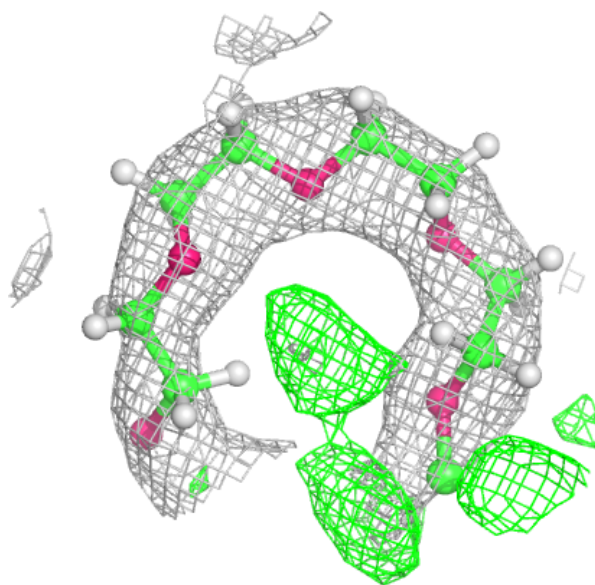
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around P6G A 1008:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.