



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 7, 2020 – 10:02 PM BST

PDB ID : 4HSI
Title : Glycoprotein B from Herpes simplex virus type 1, A504P/R505G/Q507G/N511G mutant, low-pH
Authors : Heldwein, E.E.
Deposited on : 2012-10-30
Resolution : 3.10 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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

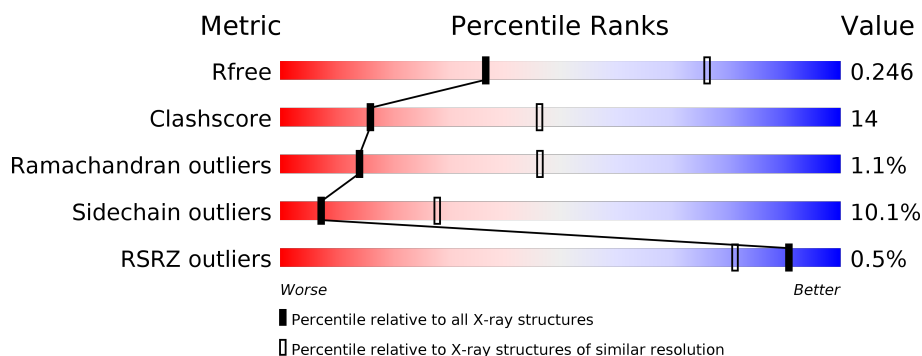
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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}	130704	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	703	<div> <div>56%</div> <div>26%</div> <div>•</div> <div>15%</div> </div>
1	B	703	<div> <div>%</div> <div>50%</div> <div>31%</div> <div>5%</div> <div>14%</div> </div>
1	C	703	<div> <div>%</div> <div>52%</div> <div>30%</div> <div>•</div> <div>14%</div> </div>
1	D	703	<div> <div>57%</div> <div>27%</div> <div>•</div> <div>14%</div> </div>
2	E	2	<div> <div>100%</div> </div>
2	F	2	<div> <div>100%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	F	1	X	-	-	-
3	NAG	C	803	X	-	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 19653 atoms, of which 0 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 Envelope glycoprotein B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	601	Total	C	N	O	S	0	0	0
			4846	3057	851	916	22			
1	B	606	Total	C	N	O	S	0	0	0
			4891	3085	862	922	22			
1	C	602	Total	C	N	O	S	0	0	0
			4856	3063	854	917	22			
1	D	605	Total	C	N	O	S	0	0	0
			4882	3079	862	919	22			

There are 156 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	28	ASP	-	expression tag	UNP P06437
A	29	PRO	-	expression tag	UNP P06437
A	30	ALA	-	expression tag	UNP P06437
A	31	ALA	-	expression tag	UNP P06437
A	32	PRO	-	expression tag	UNP P06437
A	33	THR	-	expression tag	UNP P06437
A	34	SER	-	expression tag	UNP P06437
A	35	PRO	-	expression tag	UNP P06437
A	36	GLY	-	expression tag	UNP P06437
A	37	THR	-	expression tag	UNP P06437
A	38	PRO	-	expression tag	UNP P06437
A	39	GLY	-	expression tag	UNP P06437
A	40	VAL	-	expression tag	UNP P06437
A	41	ALA	-	expression tag	UNP P06437
A	42	ALA	-	expression tag	UNP P06437
A	43	ALA	-	expression tag	UNP P06437
A	44	THR	-	expression tag	UNP P06437
A	45	GLN	-	expression tag	UNP P06437
A	46	ALA	-	expression tag	UNP P06437
A	47	ALA	-	expression tag	UNP P06437
A	48	ASN	-	expression tag	UNP P06437

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Chain	Residue	Modelled	Actual	Comment	Reference
A	49	GLY	-	expression tag	UNP P06437
A	50	GLY	-	expression tag	UNP P06437
A	51	PRO	-	expression tag	UNP P06437
A	52	ALA	-	expression tag	UNP P06437
A	53	THR	-	expression tag	UNP P06437
A	54	PRO	-	expression tag	UNP P06437
A	55	ALA	-	expression tag	UNP P06437
A	56	PRO	-	expression tag	UNP P06437
A	57	PRO	-	expression tag	UNP P06437
A	58	PRO	-	expression tag	UNP P06437
A	59	LEU	-	expression tag	UNP P06437
A	60	GLY	-	expression tag	UNP P06437
A	313	SER	THR	conflict	UNP P06437
A	443	LEU	GLN	conflict	UNP P06437
A	504	PRO	ALA	engineered mutation	UNP P06437
A	505	GLY	ARG	engineered mutation	UNP P06437
A	507	GLY	GLN	engineered mutation	UNP P06437
A	511	GLY	ASN	engineered mutation	UNP P06437
B	28	ASP	-	expression tag	UNP P06437
B	29	PRO	-	expression tag	UNP P06437
B	30	ALA	-	expression tag	UNP P06437
B	31	ALA	-	expression tag	UNP P06437
B	32	PRO	-	expression tag	UNP P06437
B	33	THR	-	expression tag	UNP P06437
B	34	SER	-	expression tag	UNP P06437
B	35	PRO	-	expression tag	UNP P06437
B	36	GLY	-	expression tag	UNP P06437
B	37	THR	-	expression tag	UNP P06437
B	38	PRO	-	expression tag	UNP P06437
B	39	GLY	-	expression tag	UNP P06437
B	40	VAL	-	expression tag	UNP P06437
B	41	ALA	-	expression tag	UNP P06437
B	42	ALA	-	expression tag	UNP P06437
B	43	ALA	-	expression tag	UNP P06437
B	44	THR	-	expression tag	UNP P06437
B	45	GLN	-	expression tag	UNP P06437
B	46	ALA	-	expression tag	UNP P06437
B	47	ALA	-	expression tag	UNP P06437
B	48	ASN	-	expression tag	UNP P06437
B	49	GLY	-	expression tag	UNP P06437
B	50	GLY	-	expression tag	UNP P06437
B	51	PRO	-	expression tag	UNP P06437

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Chain	Residue	Modelled	Actual	Comment	Reference
B	52	ALA	-	expression tag	UNP P06437
B	53	THR	-	expression tag	UNP P06437
B	54	PRO	-	expression tag	UNP P06437
B	55	ALA	-	expression tag	UNP P06437
B	56	PRO	-	expression tag	UNP P06437
B	57	PRO	-	expression tag	UNP P06437
B	58	PRO	-	expression tag	UNP P06437
B	59	LEU	-	expression tag	UNP P06437
B	60	GLY	-	expression tag	UNP P06437
B	313	SER	THR	conflict	UNP P06437
B	443	LEU	GLN	conflict	UNP P06437
B	504	PRO	ALA	engineered mutation	UNP P06437
B	505	GLY	ARG	engineered mutation	UNP P06437
B	507	GLY	GLN	engineered mutation	UNP P06437
B	511	GLY	ASN	engineered mutation	UNP P06437
C	28	ASP	-	expression tag	UNP P06437
C	29	PRO	-	expression tag	UNP P06437
C	30	ALA	-	expression tag	UNP P06437
C	31	ALA	-	expression tag	UNP P06437
C	32	PRO	-	expression tag	UNP P06437
C	33	THR	-	expression tag	UNP P06437
C	34	SER	-	expression tag	UNP P06437
C	35	PRO	-	expression tag	UNP P06437
C	36	GLY	-	expression tag	UNP P06437
C	37	THR	-	expression tag	UNP P06437
C	38	PRO	-	expression tag	UNP P06437
C	39	GLY	-	expression tag	UNP P06437
C	40	VAL	-	expression tag	UNP P06437
C	41	ALA	-	expression tag	UNP P06437
C	42	ALA	-	expression tag	UNP P06437
C	43	ALA	-	expression tag	UNP P06437
C	44	THR	-	expression tag	UNP P06437
C	45	GLN	-	expression tag	UNP P06437
C	46	ALA	-	expression tag	UNP P06437
C	47	ALA	-	expression tag	UNP P06437
C	48	ASN	-	expression tag	UNP P06437
C	49	GLY	-	expression tag	UNP P06437
C	50	GLY	-	expression tag	UNP P06437
C	51	PRO	-	expression tag	UNP P06437
C	52	ALA	-	expression tag	UNP P06437
C	53	THR	-	expression tag	UNP P06437
C	54	PRO	-	expression tag	UNP P06437

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Chain	Residue	Modelled	Actual	Comment	Reference
C	55	ALA	-	expression tag	UNP P06437
C	56	PRO	-	expression tag	UNP P06437
C	57	PRO	-	expression tag	UNP P06437
C	58	PRO	-	expression tag	UNP P06437
C	59	LEU	-	expression tag	UNP P06437
C	60	GLY	-	expression tag	UNP P06437
C	313	SER	THR	conflict	UNP P06437
C	443	LEU	GLN	conflict	UNP P06437
C	504	PRO	ALA	engineered mutation	UNP P06437
C	505	GLY	ARG	engineered mutation	UNP P06437
C	507	GLY	GLN	engineered mutation	UNP P06437
C	511	GLY	ASN	engineered mutation	UNP P06437
D	28	ASP	-	expression tag	UNP P06437
D	29	PRO	-	expression tag	UNP P06437
D	30	ALA	-	expression tag	UNP P06437
D	31	ALA	-	expression tag	UNP P06437
D	32	PRO	-	expression tag	UNP P06437
D	33	THR	-	expression tag	UNP P06437
D	34	SER	-	expression tag	UNP P06437
D	35	PRO	-	expression tag	UNP P06437
D	36	GLY	-	expression tag	UNP P06437
D	37	THR	-	expression tag	UNP P06437
D	38	PRO	-	expression tag	UNP P06437
D	39	GLY	-	expression tag	UNP P06437
D	40	VAL	-	expression tag	UNP P06437
D	41	ALA	-	expression tag	UNP P06437
D	42	ALA	-	expression tag	UNP P06437
D	43	ALA	-	expression tag	UNP P06437
D	44	THR	-	expression tag	UNP P06437
D	45	GLN	-	expression tag	UNP P06437
D	46	ALA	-	expression tag	UNP P06437
D	47	ALA	-	expression tag	UNP P06437
D	48	ASN	-	expression tag	UNP P06437
D	49	GLY	-	expression tag	UNP P06437
D	50	GLY	-	expression tag	UNP P06437
D	51	PRO	-	expression tag	UNP P06437
D	52	ALA	-	expression tag	UNP P06437
D	53	THR	-	expression tag	UNP P06437
D	54	PRO	-	expression tag	UNP P06437
D	55	ALA	-	expression tag	UNP P06437
D	56	PRO	-	expression tag	UNP P06437
D	57	PRO	-	expression tag	UNP P06437

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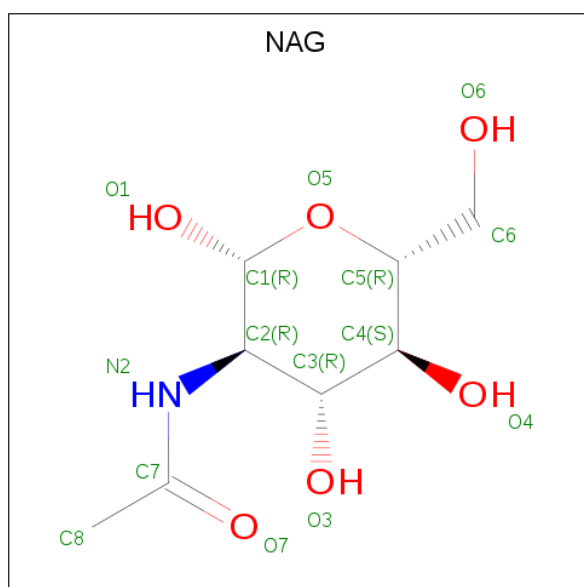
Chain	Residue	Modelled	Actual	Comment	Reference
D	58	PRO	-	expression tag	UNP P06437
D	59	LEU	-	expression tag	UNP P06437
D	60	GLY	-	expression tag	UNP P06437
D	313	SER	THR	conflict	UNP P06437
D	443	LEU	GLN	conflict	UNP P06437
D	504	PRO	ALA	engineered mutation	UNP P06437
D	505	GLY	ARG	engineered mutation	UNP P06437
D	507	GLY	GLN	engineered mutation	UNP P06437
D	511	GLY	ASN	engineered mutation	UNP P06437

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



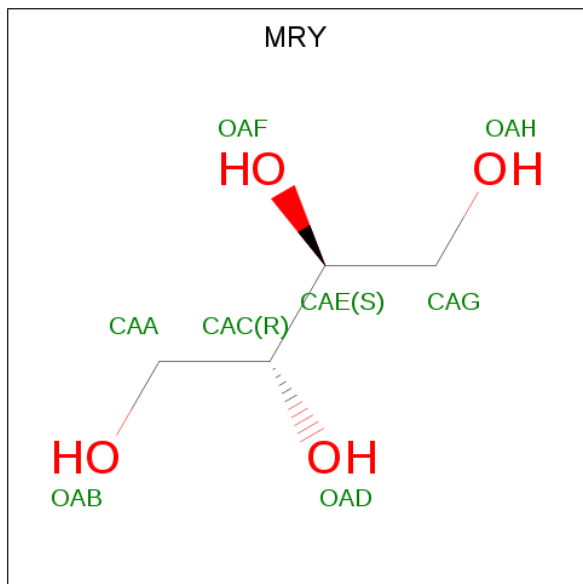
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	F	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is MESO-ERYTHRITOL (three-letter code: MRY) (formula: $C_4H_{10}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			8	4	4		
4	C	1	Total	C	O	0	0
			8	4	4		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Cl	0	0
			1	1		
5	A	1	Total	Cl	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	1	Total	Cl	0	0
			1	1		
5	C	1	Total	Cl	0	0
			1	1		

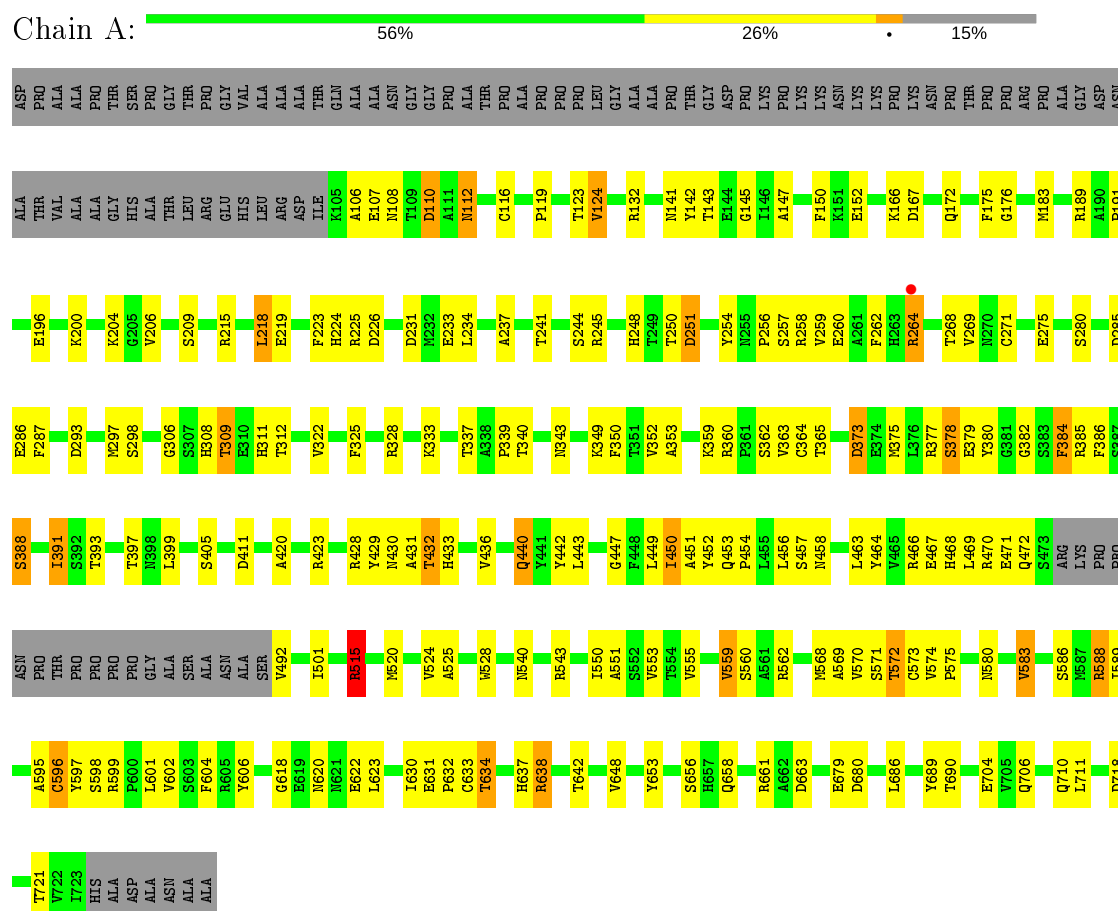
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	8	Total	O	0	0
			8	8		
6	B	1	Total	O	0	0
			1	1		
6	C	3	Total	O	0	0
			3	3		
6	D	6	Total	O	0	0
			6	6		

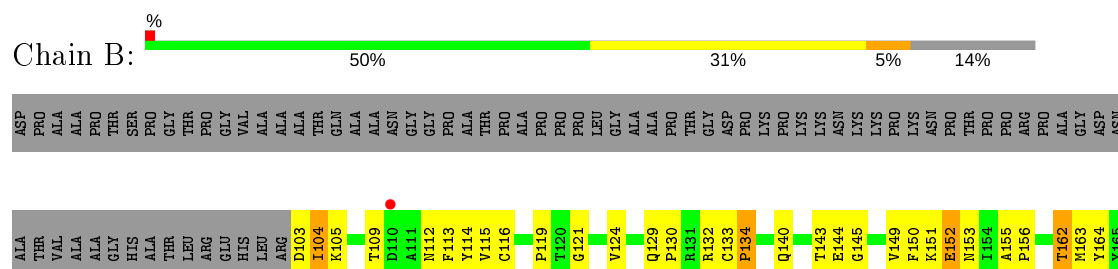
3 Residue-property plots

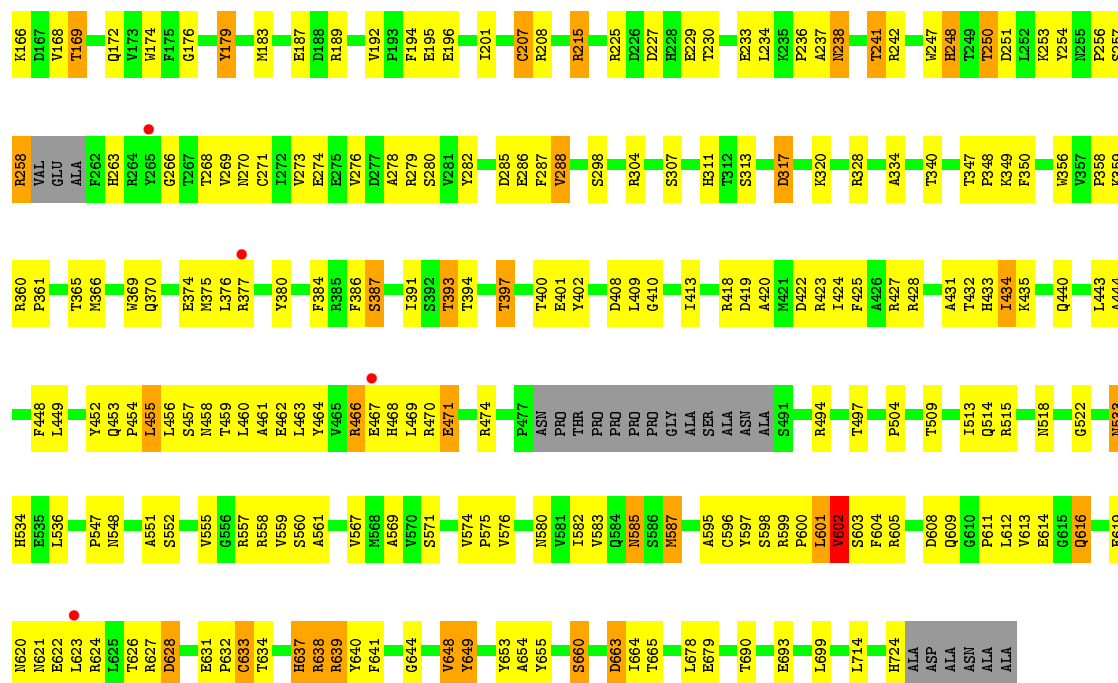
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Envelope glycoprotein B

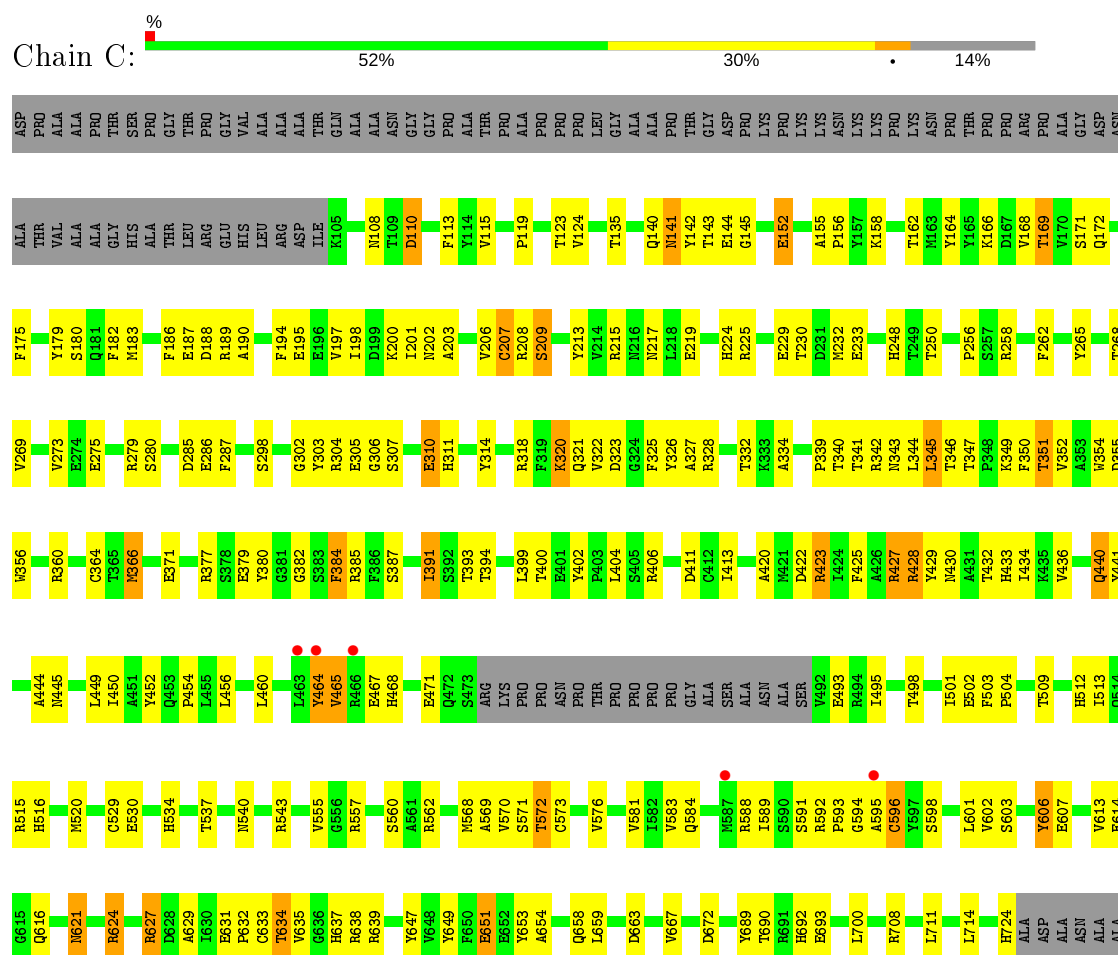


• Molecule 1: Envelope glycoprotein B



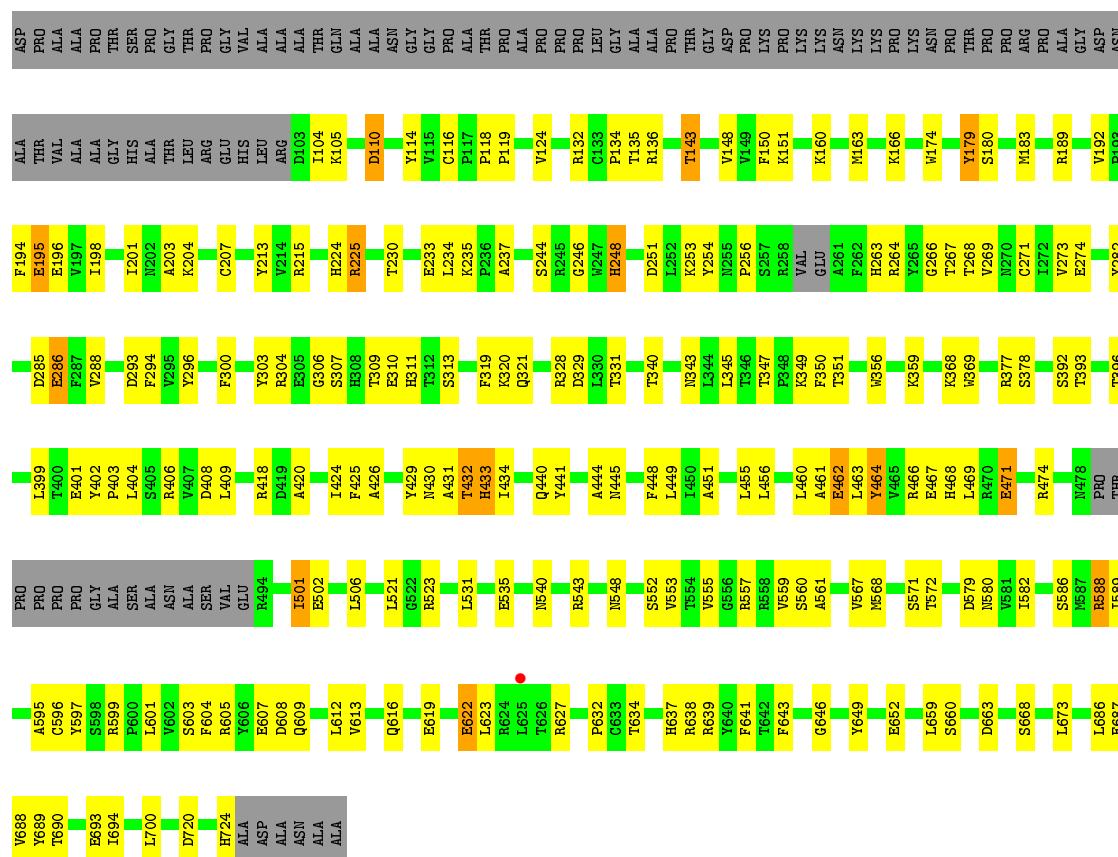


• Molecule 1: Envelope glycoprotein B



• Molecule 1: Envelope glycoprotein B

Chain D:  57% 27% 14%



- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  100%

HA61
HA62

- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  100%

HA61
HA62

4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α , β , γ	117.30Å 117.30Å 321.87Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.91 – 3.10 45.91 – 3.10	Depositor EDS
% Data completeness (in resolution range)	82.8 (45.91-3.10) 75.7 (45.91-3.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.16 (at 3.12Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.205 , 0.255 0.200 , 0.246	Depositor DCC
R_{free} test set	3774 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	52.6	Xtriage
Anisotropy	0.378	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 10.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.020 for -h,-k,l 0.459 for h,-h-k,-l 0.026 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	19653	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 73.84 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.7087e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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: CL, MRY, NAG

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	0.49	0/4966	0.64	2/6747 (0.0%)
1	B	0.45	0/5013	0.62	0/6809
1	C	0.44	0/4977	0.62	0/6762
1	D	0.46	0/5004	0.62	0/6797
All	All	0.46	0/19960	0.63	2/27115 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	515	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	A	515	ARG	NE-CZ-NH2	-5.60	117.50	120.30

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	4846	0	4657	117	1
1	B	4891	0	4703	168	0
1	C	4856	0	4664	151	1
1	D	4882	0	4694	121	0
2	E	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	28	0	25	0	0
3	A	14	0	13	0	0
3	B	28	0	26	1	0
3	C	14	0	13	0	0
3	D	28	0	26	1	0
4	A	8	0	10	0	0
4	C	8	0	10	0	0
5	A	1	0	0	1	0
5	B	1	0	0	0	0
5	C	1	0	0	1	0
5	D	1	0	0	0	0
6	A	8	0	0	0	0
6	B	1	0	0	0	0
6	C	3	0	0	0	0
6	D	6	0	0	0	0
All	All	19653	0	18866	557	2

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 14.

The worst 5 of 557 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:515:ARG:NH2	5:A:805:CL:CL	2.36	0.95
1:C:515:ARG:NH2	5:C:805:CL:CL	2.40	0.91
1:C:614:GLU:HB3	1:C:627:ARG:HH21	1.37	0.87
1:D:116:CYS:HB3	1:D:560:SER:HB3	1.63	0.81
1:B:587:MET:HB3	1:B:653:TYR:HD2	1.48	0.79

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:GLU:O	1:A:689:TYR:OH[3_555]	2.13	0.07
1:C:152:GLU:O	1:C:689:TYR:OH[3_555]	2.18	0.02

5.3 Torsion angles

5.3.1 Protein backbone

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	597/703 (85%)	529 (89%)	66 (11%)	2 (0%)	41	73
1	B	600/703 (85%)	537 (90%)	54 (9%)	9 (2%)	10	39
1	C	598/703 (85%)	527 (88%)	64 (11%)	7 (1%)	13	44
1	D	599/703 (85%)	537 (90%)	54 (9%)	8 (1%)	12	42
All	All	2394/2812 (85%)	2130 (89%)	238 (10%)	26 (1%)	14	46

5 of 26 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	237	ALA
1	B	413	ILE
1	C	413	ILE
1	D	462	GLU
1	C	465	VAL

5.3.2 Protein sidechains

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	522/592 (88%)	463 (89%)	59 (11%)	6	23
1	B	528/592 (89%)	466 (88%)	62 (12%)	5	22
1	C	523/592 (88%)	470 (90%)	53 (10%)	7	28
1	D	526/592 (89%)	489 (93%)	37 (7%)	15	45
All	All	2099/2368 (89%)	1888 (90%)	211 (10%)	7	28

5 of 211 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	459	THR
1	B	663	ASP
1	D	432	THR
1	B	471	GLU
1	B	616	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	440	GLN
1	A	658	GLN
1	C	468	HIS
1	D	440	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

4 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	$\# Z > 2$
2	NAG	E	1	1,2	14,14,15	0.63	0	17,19,21	1.19	2 (11%)
2	NAG	E	2	2	14,14,15	0.56	0	17,19,21	1.08	2 (11%)
2	NAG	F	1	1,2	14,14,15	0.61	0	17,19,21	1.13	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	F	2	2	14,14,15	0.56	0	17,19,21	1.82	4 (23%)

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
2	NAG	E	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	3/6/23/26	0/1/1/1
2	NAG	F	1	1,2	1/1/5/7	0/6/23/26	0/1/1/1
2	NAG	F	2	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	2	NAG	O5-C1-C2	-4.60	104.02	111.29
2	F	2	NAG	C3-C4-C5	3.79	117.00	110.24
2	E	1	NAG	C4-C3-C2	3.40	116.00	111.02
2	F	1	NAG	C4-C3-C2	2.61	114.84	111.02
2	F	2	NAG	O5-C5-C4	2.46	116.81	110.83

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	F	1	NAG	C1

5 of 7 torsion outliers are listed below:

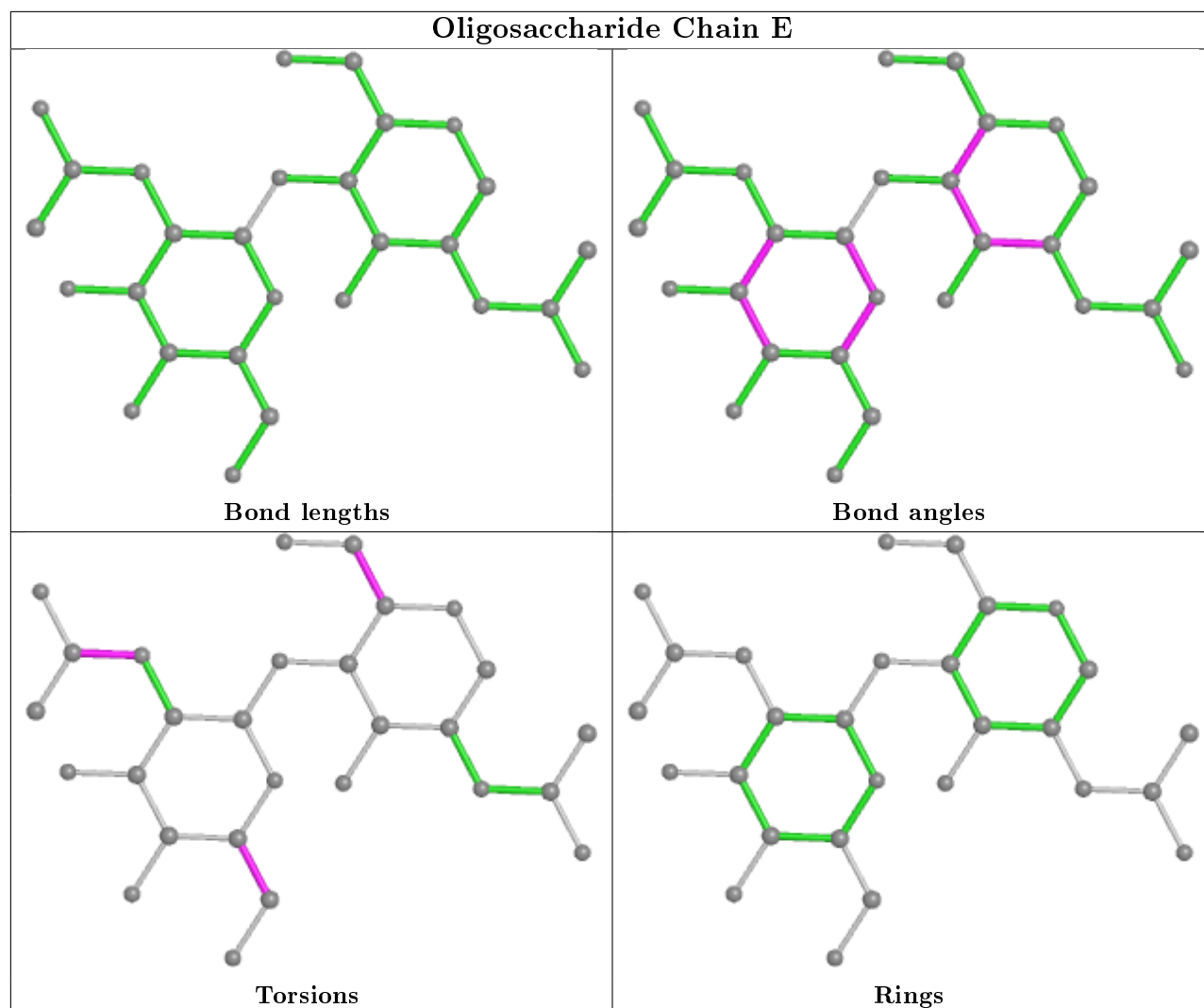
Mol	Chain	Res	Type	Atoms
2	F	2	NAG	C8-C7-N2-C2
2	F	2	NAG	O7-C7-N2-C2
2	E	1	NAG	O5-C5-C6-O6
2	E	2	NAG	C8-C7-N2-C2
2	E	2	NAG	O7-C7-N2-C2

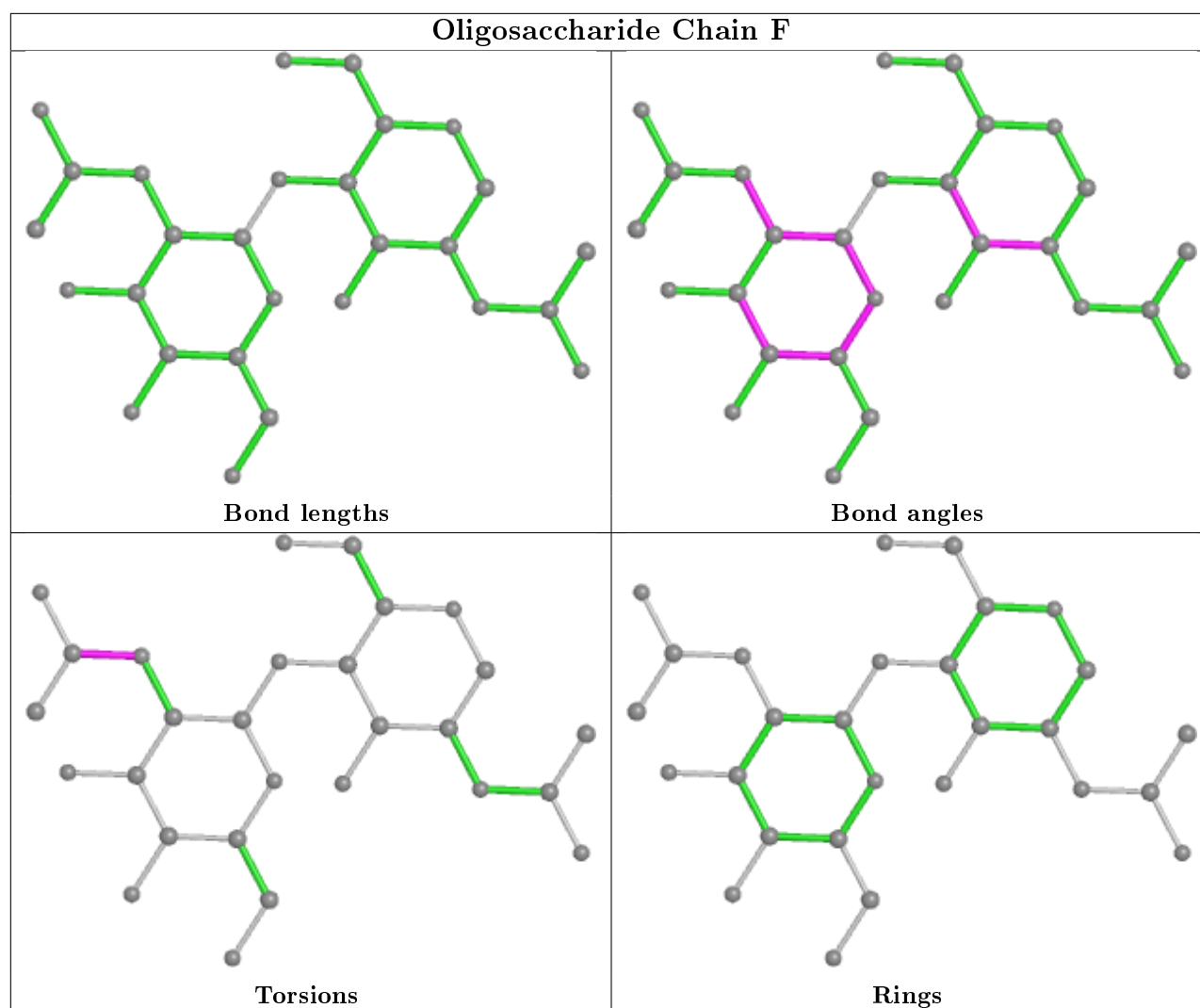
There are no ring outliers.

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 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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	$\# Z > 2$
3	NAG	D	802	1	14,14,15	0.60	0	17,19,21	1.33	2 (11%)
3	NAG	C	803	1	14,14,15	0.46	0	17,19,21	1.67	2 (11%)
3	NAG	D	801	1	14,14,15	0.43	0	17,19,21	1.50	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	B	801	1	14,14,15	0.55	0	17,19,21	1.77	4 (23%)
3	NAG	A	803	1	14,14,15	0.73	0	17,19,21	1.16	2 (11%)
4	MRY	C	804	-	7,7,7	0.33	0	8,8,8	1.07	0
4	MRY	A	804	-	7,7,7	0.39	0	8,8,8	0.87	0
3	NAG	B	802	1	14,14,15	0.59	0	17,19,21	1.37	1 (5%)

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	D	802	1	-	2/6/23/26	0/1/1/1
3	NAG	C	803	1	1/1/5/7	4/6/23/26	0/1/1/1
3	NAG	D	801	1	-	4/6/23/26	0/1/1/1
3	NAG	B	801	1	-	0/6/23/26	0/1/1/1
3	NAG	A	803	1	-	3/6/23/26	0/1/1/1
4	MRY	C	804	-	-	0/8/8/8	-
4	MRY	A	804	-	-	0/8/8/8	-
3	NAG	B	802	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	803	NAG	O5-C1-C2	-5.40	102.77	111.29
3	D	801	NAG	C1-O5-C5	4.43	118.20	112.19
3	B	801	NAG	C1-O5-C5	4.18	117.85	112.19
3	B	802	NAG	C2-N2-C7	-3.70	117.63	122.90
3	B	801	NAG	C2-N2-C7	3.69	128.16	122.90

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	C	803	NAG	C1

5 of 15 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	802	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
3	C	803	NAG	C8-C7-N2-C2
3	C	803	NAG	O7-C7-N2-C2
3	D	801	NAG	C8-C7-N2-C2
3	D	801	NAG	O7-C7-N2-C2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	801	NAG	1	0
3	B	801	NAG	1	0

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, 95th 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 ²)	Q<0.9
1	A	601/703 (85%)	-0.18	1 (0%) 95 90	1, 37, 84, 140	0
1	B	606/703 (86%)	-0.13	5 (0%) 86 72	3, 44, 95, 185	0
1	C	602/703 (85%)	-0.14	5 (0%) 86 72	7, 44, 98, 157	0
1	D	605/703 (86%)	-0.18	1 (0%) 95 90	2, 39, 78, 147	1 (0%)
All	All	2414/2812 (85%)	-0.16	12 (0%) 91 81	1, 41, 89, 185	1 (0%)

The worst 5 of 12 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	265	TYR	3.0
1	B	467	GLU	2.8
1	C	464	TYR	2.8
1	A	264	ARG	2.7
1	B	623	LEU	2.7

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, 95th 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.

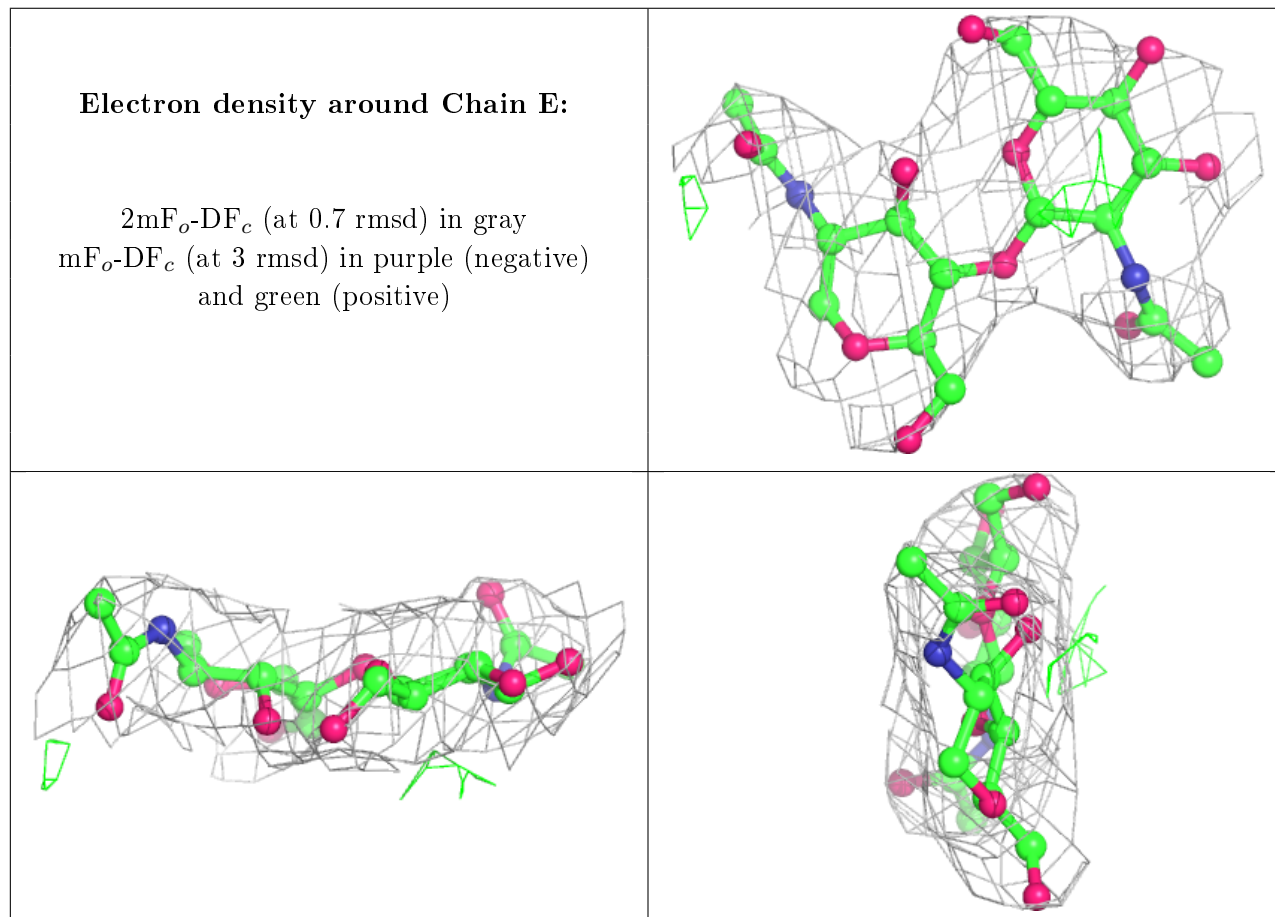
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	F	1	14/15	0.90	0.19	88,88,88,88	0
2	NAG	F	2	14/15	0.90	0.17	92,92,92,92	0

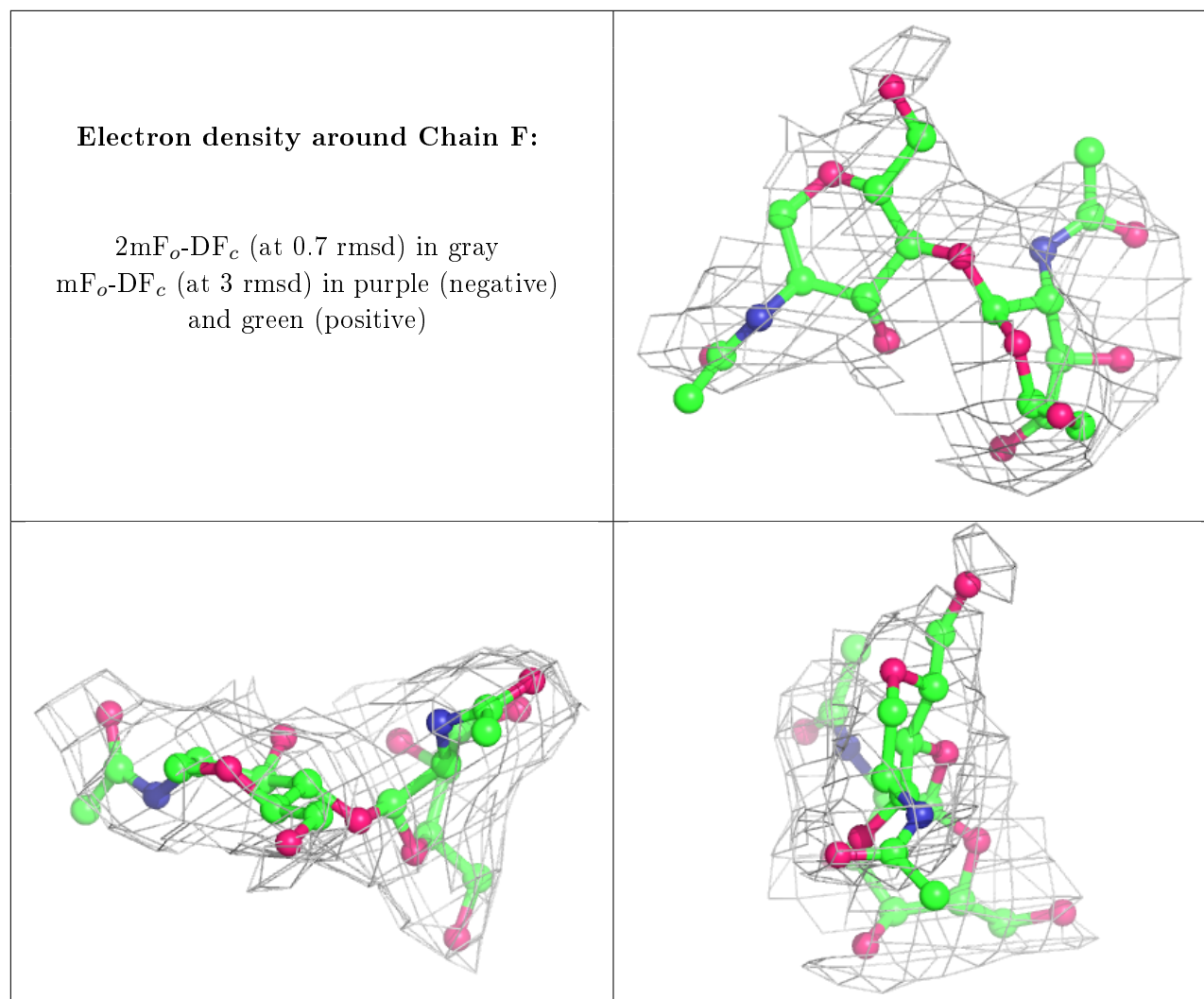
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAG	E	2	14/15	0.92	0.16	86,86,86,86	0
2	NAG	E	1	14/15	0.93	0.19	92,92,92,92	0

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 ⓘ

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, 95th 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(\AA^2)	Q<0.9
5	CL	B	803	1/1	0.81	0.20	67,67,67,67	0
3	NAG	A	803	14/15	0.88	0.15	71,71,71,71	0
3	NAG	C	803	14/15	0.88	0.15	71,71,71,71	0
3	NAG	B	801	14/15	0.91	0.19	69,69,69,69	0
4	MRY	C	804	8/8	0.94	0.23	34,34,34,34	0
4	MRY	A	804	8/8	0.94	0.23	44,44,44,44	0
5	CL	D	803	1/1	0.94	0.26	71,71,71,71	0
3	NAG	D	801	14/15	0.94	0.15	71,71,71,71	0
5	CL	A	805	1/1	0.96	0.13	30,30,30,30	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAG	B	802	14/15	0.96	0.14	31,31,31,31	0
5	CL	C	805	1/1	0.97	0.12	32,32,32,32	0
3	NAG	D	802	14/15	0.97	0.14	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.