



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 10, 2020 – 10:20 AM BST

PDB ID : 6UYD
Title : Structure of Hepatitis C Virus Envelope Glycoprotein E2mc3-v1 redesigned core from genotype 1a bound to broadly neutralizing antibody AR3C
Authors : Tzarum, N.; Wilson, I.A.; Zhu, J.
Deposited on : 2019-11-13
Resolution : 1.90 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

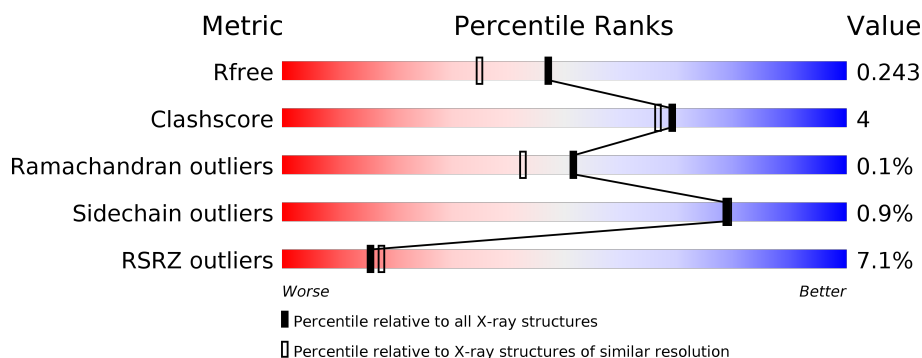
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	E	177	<div> <div>10%</div> <div> <div>68%</div> <div>11%</div> <div>19%</div> </div> </div>
1	F	177	<div> <div>5%</div> <div> <div>76%</div> <div>12%</div> <div>11%</div> </div> </div>
2	A	233	<div> <div>10%</div> <div> <div>85%</div> <div>10%</div> <div>5%</div> </div> </div>
2	H	233	<div> <div>6%</div> <div> <div>89%</div> <div>6%</div> <div>5%</div> </div> </div>
3	B	214	<div> <div>8%</div> <div> <div>93%</div> <div>6%</div> </div> </div>
3	L	214	<div> <div>%</div> <div> <div>93%</div> <div>5%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
4	C	2	 50%50%
4	D	2	 50%50%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 9525 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 E2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	143	Total	C	N	O	S	0	0	0
			1118	719	189	197	13			
1	F	158	Total	C	N	O	S	0	0	0
			1230	782	210	225	13			

- Molecule 2 is a protein called Fab AR3C heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	222	Total	C	N	O	S	0	0	0
			1659	1044	281	327	7			
2	A	222	Total	C	N	O	S	0	0	0
			1656	1042	280	326	8			

- Molecule 3 is a protein called Fab AR3C light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	211	Total	C	N	O	S	0	0	0
			1612	1005	277	326	4			
3	B	212	Total	C	N	O	S	0	0	0
			1618	1008	278	327	5			

- Molecule 4 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
4	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

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

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	E	1	Total	C	N	O	0	0
			14	8	1	5		
5	E	1	Total	C	N	O	0	0
			14	8	1	5		
5	E	1	Total	C	N	O	0	0
			14	8	1	5		
5	F	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	E	38	Total	O	0	0
			38	38		
6	H	134	Total	O	0	0
			134	134		
6	L	104	Total	O	0	0
			104	104		
6	A	100	Total	O	0	0
			100	100		

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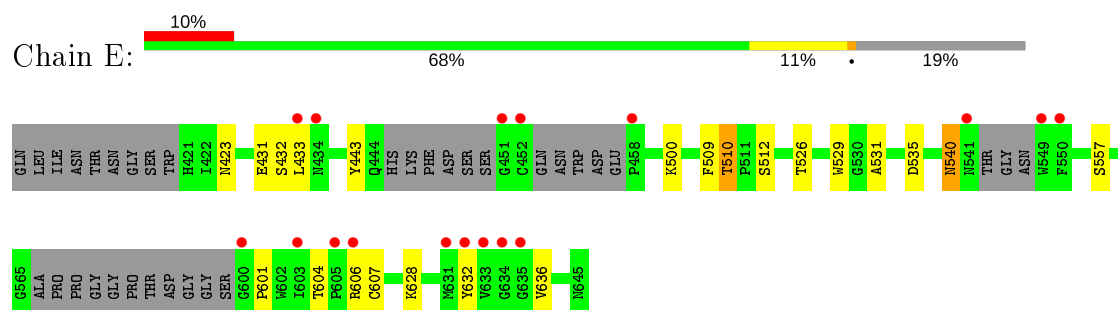
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	57	Total	O	0	0
			57	57		
6	F	87	Total	O	0	0
			87	87		

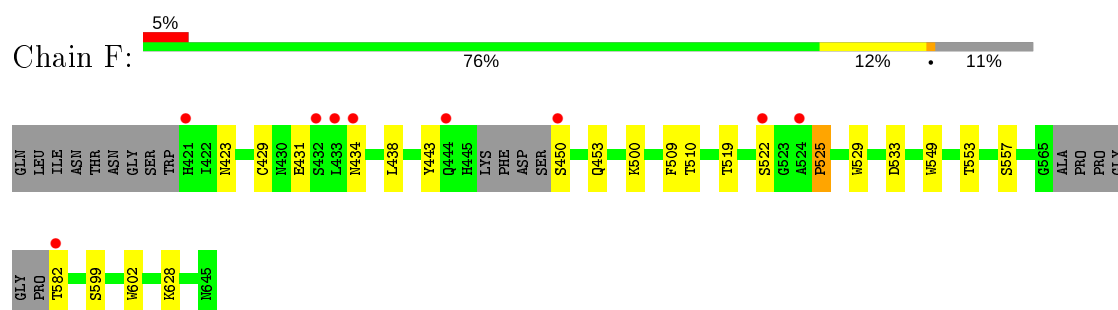
3 Residue-property plots [i](#)

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.

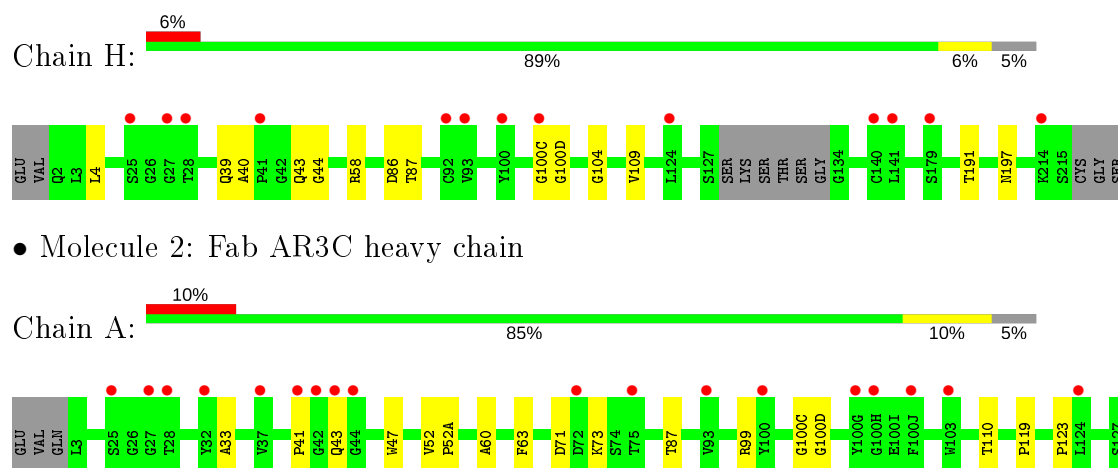
- Molecule 1: Envelope glycoprotein E2



- Molecule 1: Envelope glycoprotein E2



- Molecule 2: Fab AR3C heavy chain



- Molecule 2: Fab AR3C heavy chain



- Molecule 3: Fab AR3C light chain



- Molecule 3: Fab AR3C light chain



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	88.49Å 93.66Å 96.80Å 90.00° 100.48° 90.00°	Depositor
Resolution (Å)	29.91 – 1.90 29.91 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.8 (29.91-1.90) 98.8 (29.91-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.16 (at 1.89Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.208 , 0.243 0.208 , 0.243	Depositor DCC
R_{free} test set	6108 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	35.2	Xtriage
Anisotropy	0.517	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9525	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.23% of the height of the origin peak. No significant pseudotranslation is detected.*

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: 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	E	0.47	0/1156	0.61	0/1580
1	F	0.45	0/1273	0.60	0/1744
2	A	0.44	0/1695	0.58	0/2306
2	H	0.48	0/1698	0.61	0/2310
3	B	0.35	0/1652	0.53	0/2243
3	L	0.44	0/1646	0.60	0/2235
All	All	0.44	0/9120	0.59	0/12418

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	E	1118	0	1037	14	0
1	F	1230	0	1124	15	0
2	A	1656	0	1610	18	0
2	H	1659	0	1614	8	0
3	B	1618	0	1574	9	0
3	L	1612	0	1570	5	0
4	C	28	0	25	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	28	0	25	1	0
5	E	42	0	39	1	0
5	F	14	0	13	0	0
6	A	100	0	0	0	0
6	B	57	0	0	1	0
6	E	38	0	0	1	0
6	F	87	0	0	3	0
6	H	134	0	0	2	0
6	L	104	0	0	0	0
All	All	9525	0	8631	65	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 4.

The worst 5 of 65 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)
2:A:41:PRO:O	2:A:43:GLN:NE2	2.03	0.92
1:F:429:CYS:HB3	1:F:438:LEU:HD21	1.67	0.75
1:E:604:THR:HG22	1:E:606:ARG:H	1.52	0.74
1:F:450:SER:HB3	1:F:453:GLN:HB2	1.70	0.73
3:L:142:ARG:NH2	3:L:163:VAL:HG11	2.09	0.67

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	E	133/177 (75%)	129 (97%)	3 (2%)	1 (1%)	19	9
1	F	152/177 (86%)	148 (97%)	4 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	A	218/233 (94%)	212 (97%)	6 (3%)	0	100	100
2	H	218/233 (94%)	209 (96%)	9 (4%)	0	100	100
3	B	210/214 (98%)	203 (97%)	7 (3%)	0	100	100
3	L	209/214 (98%)	203 (97%)	6 (3%)	0	100	100
All	All	1140/1248 (91%)	1104 (97%)	35 (3%)	1 (0%)	51	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	510	THR

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	E	122/149 (82%)	118 (97%)	4 (3%)	38	29
1	F	134/149 (90%)	131 (98%)	3 (2%)	52	47
2	A	186/195 (95%)	186 (100%)	0	100	100
2	H	186/195 (95%)	185 (100%)	1 (0%)	88	89
3	B	182/184 (99%)	182 (100%)	0	100	100
3	L	181/184 (98%)	180 (99%)	1 (1%)	86	87
All	All	991/1056 (94%)	982 (99%)	9 (1%)	78	79

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

Mol	Chain	Res	Type
2	H	197	ASN
1	F	525	PRO
1	F	443	TYR
1	E	540	ASN
3	L	33	LEU

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

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
4	NAG	C	1	1,4	14,14,15	0.18	0	17,19,21	0.61	1 (5%)
4	NAG	C	2	4	14,14,15	0.26	0	17,19,21	0.65	1 (5%)
4	NAG	D	1	1,4	14,14,15	0.46	0	17,19,21	0.37	0
4	NAG	D	2	4	14,14,15	0.30	0	17,19,21	0.62	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	NAG	C	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	C	2	4	-	0/6/23/26	0/1/1/1
4	NAG	D	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	D	2	4	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	2	NAG	C1-O5-C5	2.27	115.27	112.19
4	C	1	NAG	C1-O5-C5	2.05	114.97	112.19

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	2	NAG	C4-C5-C6-O6
4	D	2	NAG	O5-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	1	NAG	1	0
4	C	1	NAG	1	0

5.6 Ligand geometry

4 ligands 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	$\# Z > 2$
5	NAG	E	703	1	14,14,15	0.35	0	17,19,21	0.57	0
5	NAG	F	701	1	14,14,15	0.44	0	17,19,21	0.56	0
5	NAG	E	702	1	14,14,15	0.42	0	17,19,21	1.26	1 (5%)
5	NAG	E	701	1	14,14,15	0.60	0	17,19,21	0.66	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
5	NAG	E	703	1	-	2/6/23/26	0/1/1/1
5	NAG	F	701	1	-	0/6/23/26	0/1/1/1
5	NAG	E	702	1	-	3/6/23/26	0/1/1/1
5	NAG	E	701	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	702	NAG	C2-N2-C7	4.45	129.24	122.90
5	E	701	NAG	C1-O5-C5	2.07	115.00	112.19

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	701	NAG	O5-C5-C6-O6
5	E	703	NAG	O5-C5-C6-O6
5	E	703	NAG	C4-C5-C6-O6
5	E	701	NAG	C4-C5-C6-O6
5	E	702	NAG	C8-C7-N2-C2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	702	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	E	143/177 (80%)	0.58	17 (11%) 4 5	27, 44, 78, 98	0
1	F	158/177 (89%)	0.23	9 (5%) 23 26	26, 42, 72, 99	0
2	A	222/233 (95%)	0.40	23 (10%) 6 7	30, 45, 75, 104	0
2	H	222/233 (95%)	0.24	13 (5%) 22 25	25, 40, 70, 82	0
3	B	212/214 (99%)	0.53	18 (8%) 10 12	38, 54, 75, 97	0
3	L	211/214 (98%)	0.04	3 (1%) 75 77	27, 42, 61, 76	0
All	All	1168/1248 (93%)	0.33	83 (7%) 16 17	25, 44, 73, 104	0

The worst 5 of 83 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	25	SER	6.6
2	H	27	GLY	6.4
1	E	433	LEU	6.2
1	E	451	GLY	6.1
1	E	632	TYR	5.3

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(\AA^2)	Q<0.9
4	NAG	C	2	14/15	0.85	0.28	64,80,91,96	0
4	NAG	D	2	14/15	0.92	0.22	62,74,96,98	0
4	NAG	D	1	14/15	0.94	0.10	45,57,69,71	0
4	NAG	C	1	14/15	0.95	0.12	45,54,60,71	0

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, 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	NAG	E	702	14/15	0.57	0.33	67,88,97,100	0
5	NAG	E	703	14/15	0.69	0.36	91,100,105,106	0
5	NAG	F	701	14/15	0.79	0.37	82,90,101,108	0
5	NAG	E	701	14/15	0.80	0.31	82,93,101,117	0

6.5 Other polymers [i](#)

There are no such residues in this entry.