



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 19, 2016 – 10:56 PM GMT

PDB ID : 5EJB
Title : Crystal structure of prefusion Hendra virus F protein
Authors : Wong, J.W.; Jardetzky, T.S.; Paterson, R.G.; Lamb, R.A.
Deposited on : 2015-11-01
Resolution : 3.20 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
<http://wwpdb.org/validation/2016/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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20026982

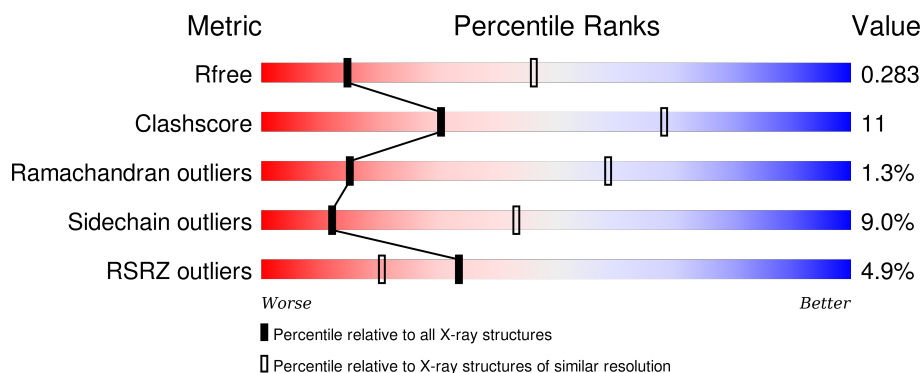
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	91344	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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. 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	501	<div> <div>3%</div> <div>62% 23% • 12%</div> </div>
1	B	501	<div> <div>4%</div> <div>63% 23% • 11%</div> </div>
1	C	501	<div> <div>7%</div> <div>66% 21% • 9%</div> </div>
1	D	501	<div> <div>5%</div> <div>64% 23% • 8%</div> </div>
1	E	501	<div> <div>5%</div> <div>64% 21% • 13%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	501	

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	A	602	-	-	-	X

2 Entry composition [i](#)

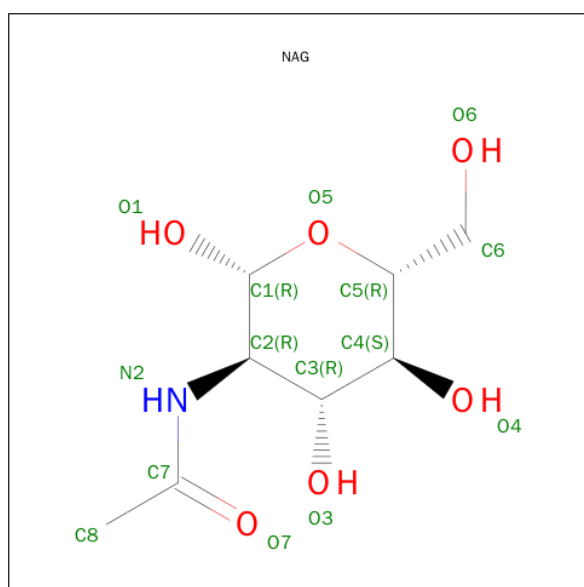
There are 3 unique types of molecules in this entry. The entry contains 20415 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Fusion glycoprotein F0.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	447	Total	C	N	O	S	0	0	0
			3376	2135	553	670	18			
1	C	454	Total	C	N	O	S	0	0	0
			3390	2142	558	672	18			
1	D	460	Total	C	N	O	S	0	0	0
			3441	2174	566	683	18			
1	F	436	Total	C	N	O	S	0	0	0
			3298	2086	540	654	18			
1	E	436	Total	C	N	O	S	0	0	0
			3298	2086	540	654	18			
1	A	440	Total	C	N	O	S	0	0	0
			3330	2107	545	660	18			

- Molecule 2 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).

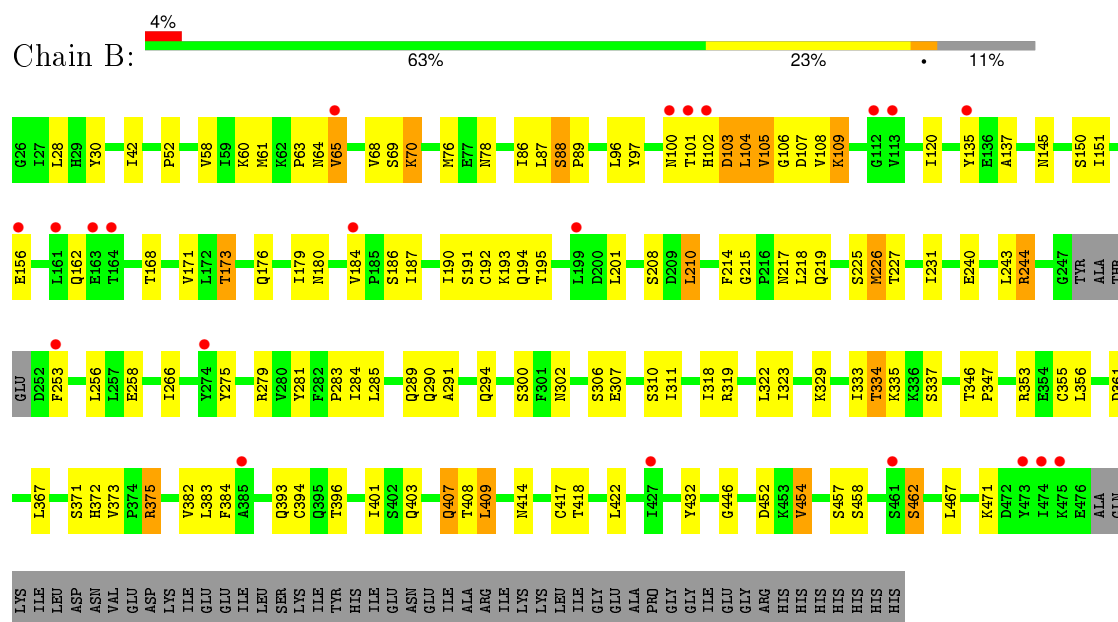


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

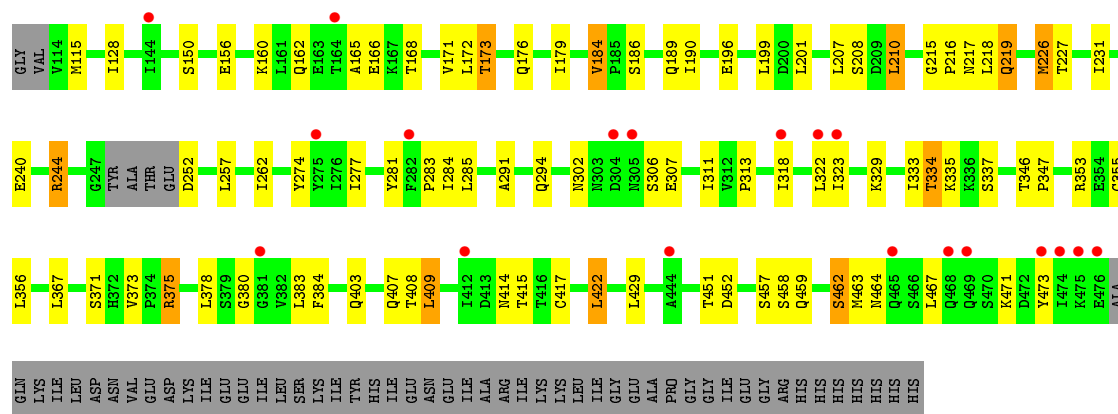
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ($\text{RSRZ} > 2$). 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.

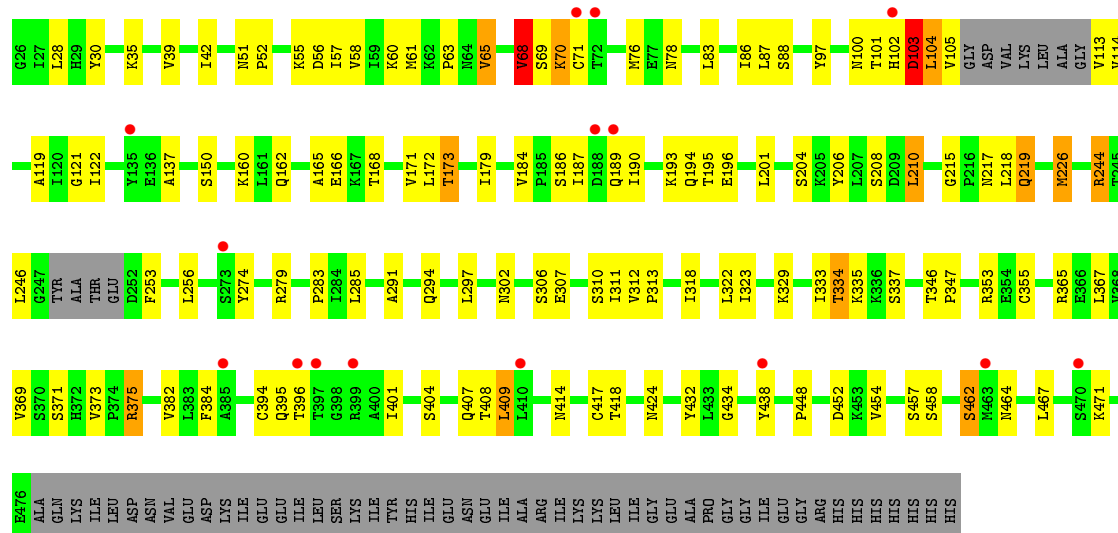
• Molecule 1: Fusion glycoprotein F0







● Molecule 1: Fusion glycoprotein F0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	108.61Å 163.50Å 147.94Å 90.00° 94.13° 90.00°	Depositor
Resolution (Å)	49.62 – 3.20 49.62 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.62-3.20) 99.7 (49.62-3.20)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 3.19Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.253 , 0.283 0.253 , 0.283	Depositor DCC
R_{free} test set	4233 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	75.1	Xtriage
Anisotropy	0.888	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 60.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 84644 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	20415	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.87% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, SO4

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	0.54	0/3375	0.73	1/4593 (0.0%)
1	B	0.59	0/3422	0.75	0/4657
1	C	0.53	0/3435	0.71	0/4677
1	D	0.59	0/3487	0.75	1/4748 (0.0%)
1	E	0.52	0/3343	0.71	0/4548
1	F	0.54	0/3343	0.72	0/4548
All	All	0.55	0/20405	0.73	2/27771 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	D	0	1
All	All	0	3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	319	ARG	NE-CZ-NH2	-5.66	117.47	120.30
1	A	253	PHE	N-CA-CB	-5.20	101.25	110.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	103	ASP	Peptide
1	B	102	HIS	Peptide
1	D	102	HIS	Peptide

5.2 Too-close contacts

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	3330	0	3324	87	0
1	B	3376	0	3377	86	1
1	C	3390	0	3336	75	0
1	D	3441	0	3403	88	0
1	E	3298	0	3288	75	0
1	F	3298	0	3290	84	1
2	A	28	0	26	0	0
2	B	28	0	25	0	0
2	C	42	0	38	0	0
2	D	70	0	64	2	0
2	E	56	0	52	1	0
2	F	28	0	26	2	0
3	A	5	0	0	0	0
3	B	5	0	0	1	0
3	C	5	0	0	0	0
3	D	5	0	0	1	0
3	E	5	0	0	0	0
3	F	5	0	0	1	0
All	All	20415	0	20249	449	1

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 11.

The worst 5 of 449 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:F:318:ILE:HG12	1:F:323:ILE:HG22	1.49	0.92
1:E:318:ILE:HG12	1:E:323:ILE:HG22	1.52	0.91
1:C:65:VAL:HG21	1:C:76:MET:CE	2.00	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:307:GLU:HG3	1:D:375:ARG:HH21	1.36	0.91
1:E:65:VAL:HG21	1:E:76:MET:CE	2.01	0.91

All (1) 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:B:393:GLN:OE1	1:F:289:GLN:OE1[2_656]	2.12	0.08

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	434/501 (87%)	393 (91%)	34 (8%)	7 (2%)	12	54
1	B	443/501 (88%)	392 (88%)	43 (10%)	8 (2%)	11	51
1	C	448/501 (89%)	399 (89%)	42 (9%)	7 (2%)	12	54
1	D	456/501 (91%)	406 (89%)	41 (9%)	9 (2%)	9	48
1	E	430/501 (86%)	388 (90%)	40 (9%)	2 (0%)	34	78
1	F	430/501 (86%)	386 (90%)	42 (10%)	2 (0%)	34	78
All	All	2641/3006 (88%)	2364 (90%)	242 (9%)	35 (1%)	15	59

5 of 35 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	69	SER
1	B	104	LEU
1	B	107	ASP
1	B	371	SER
1	C	27	ILE

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	377/443 (85%)	341 (90%)	36 (10%)	11	40
1	B	382/443 (86%)	347 (91%)	35 (9%)	11	41
1	C	374/443 (84%)	342 (91%)	32 (9%)	13	46
1	D	382/443 (86%)	345 (90%)	37 (10%)	10	39
1	E	373/443 (84%)	342 (92%)	31 (8%)	14	49
1	F	373/443 (84%)	340 (91%)	33 (9%)	12	45
All	All	2261/2658 (85%)	2057 (91%)	204 (9%)	12	43

5 of 204 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	334	THR
1	F	201	LEU
1	A	306	SER
1	D	367	LEU
1	D	462	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 33 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	290	GLN
1	F	162	GLN
1	A	314	ASN
1	D	314	ASN
1	D	325	ASN

5.3.3 RNA [i](#)

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

24 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	$\# Z > 2$
2	NAG	A	601	1	14,14,15	0.67	0	15,19,21	1.73	2 (13%)
2	NAG	A	602	1	14,14,15	0.54	0	15,19,21	2.31	3 (20%)
3	SO4	A	603	-	4,4,4	0.74	0	6,6,6	1.23	1 (16%)
2	NAG	B	601	1,2	14,14,15	0.58	0	15,19,21	1.62	2 (13%)
2	NAG	B	602	2	14,14,15	0.47	0	15,19,21	1.16	1 (6%)
3	SO4	B	603	-	4,4,4	0.38	0	6,6,6	0.19	0
2	NAG	C	601	1,2	14,14,15	0.82	0	15,19,21	1.66	1 (6%)
2	NAG	C	602	2	14,14,15	0.68	0	15,19,21	1.47	3 (20%)
2	NAG	C	603	1	14,14,15	0.81	1 (7%)	15,19,21	2.12	3 (20%)
3	SO4	C	604	-	4,4,4	1.12	0	6,6,6	2.01	1 (16%)
2	NAG	D	601	1	14,14,15	0.75	0	15,19,21	1.58	3 (20%)
2	NAG	D	602	1	14,14,15	0.42	0	15,19,21	1.54	3 (20%)
2	NAG	D	603	1,2	14,14,15	0.81	0	15,19,21	2.23	7 (46%)
2	NAG	D	604	2	14,14,15	0.55	0	15,19,21	1.53	3 (20%)
2	NAG	D	605	1	14,14,15	0.58	0	15,19,21	1.15	1 (6%)
3	SO4	D	606	-	4,4,4	0.43	0	6,6,6	0.47	0
2	NAG	E	601	1	14,14,15	0.61	0	15,19,21	1.88	3 (20%)
2	NAG	E	602	1	14,14,15	0.55	0	15,19,21	1.58	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	E	603	1	14,14,15	0.52	0	15,19,21	1.87	3 (20%)
2	NAG	E	604	1	14,14,15	0.58	0	15,19,21	1.23	3 (20%)
3	SO4	E	605	-	4,4,4	0.64	0	6,6,6	0.96	0
2	NAG	F	601	1	14,14,15	0.62	0	15,19,21	1.49	2 (13%)
2	NAG	F	602	1	14,14,15	0.63	0	15,19,21	1.39	3 (20%)
3	SO4	F	603	-	4,4,4	0.77	0	6,6,6	0.70	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
2	NAG	A	601	1	-	0/6/23/26	0/1/1/1
2	NAG	A	602	1	-	0/6/23/26	0/1/1/1
3	SO4	A	603	-	-	0/0/0/0	0/0/0/0
2	NAG	B	601	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	602	2	-	0/6/23/26	0/1/1/1
3	SO4	B	603	-	-	0/0/0/0	0/0/0/0
2	NAG	C	601	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	602	2	-	0/6/23/26	0/1/1/1
2	NAG	C	603	1	-	0/6/23/26	0/1/1/1
3	SO4	C	604	-	-	0/0/0/0	0/0/0/0
2	NAG	D	601	1	-	0/6/23/26	0/1/1/1
2	NAG	D	602	1	-	0/6/23/26	0/1/1/1
2	NAG	D	603	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	604	2	-	0/6/23/26	0/1/1/1
2	NAG	D	605	1	-	0/6/23/26	0/1/1/1
3	SO4	D	606	-	-	0/0/0/0	0/0/0/0
2	NAG	E	601	1	-	0/6/23/26	0/1/1/1
2	NAG	E	602	1	-	0/6/23/26	0/1/1/1
2	NAG	E	603	1	-	0/6/23/26	0/1/1/1
2	NAG	E	604	1	-	0/6/23/26	0/1/1/1
3	SO4	E	605	-	-	0/0/0/0	0/0/0/0
2	NAG	F	601	1	-	0/6/23/26	0/1/1/1
2	NAG	F	602	1	-	0/6/23/26	0/1/1/1
3	SO4	F	603	-	-	0/0/0/0	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	603	NAG	C1-C2	2.15	1.55	1.52

The worst 5 of 50 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	602	NAG	C4-C3-C2	-4.11	104.97	111.34
2	B	601	NAG	C2-N2-C7	-3.87	118.08	123.11
2	D	603	NAG	O3-C3-C4	-2.25	105.27	110.36
2	F	602	NAG	O7-C7-C8	-2.19	118.03	122.07
2	D	602	NAG	C2-N2-C7	-2.07	120.41	123.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	603	SO4	1	0
2	D	605	NAG	2	0
3	D	606	SO4	1	0
2	E	603	NAG	1	0
2	F	602	NAG	2	0
3	F	603	SO4	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	440/501 (87%)	0.27	15 (3%)	49	34	43, 91, 149, 193	0
1	B	447/501 (89%)	0.30	21 (4%)	35	22	52, 81, 147, 170	0
1	C	454/501 (90%)	0.43	35 (7%)	16	9	51, 94, 151, 183	0
1	D	460/501 (91%)	0.33	23 (5%)	32	19	54, 84, 155, 194	0
1	E	436/501 (87%)	0.35	23 (5%)	30	17	55, 91, 154, 203	0
1	F	436/501 (87%)	0.22	15 (3%)	49	34	48, 90, 149, 184	0
All	All	2673/3006 (88%)	0.32	132 (4%)	33	20	43, 89, 151, 203	0

The worst 5 of 132 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	164	THR	7.1
1	F	102	HIS	7.1
1	A	102	HIS	6.7
1	B	164	THR	5.8
1	D	474	ILE	4.8

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](#)

There are no carbohydrates in this entry.

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	NAG	A	602	14/15	0.61	0.41	4.80	108,123,132,139	0
3	SO4	D	606	5/5	0.88	0.29	1.71	111,124,131,137	0
2	NAG	F	602	14/15	0.75	0.39	1.61	98,112,127,129	0
2	NAG	D	603	14/15	0.76	0.29	1.37	66,92,103,111	0
2	NAG	E	603	14/15	0.60	0.33	0.47	110,134,146,149	0
3	SO4	B	603	5/5	0.82	0.22	0.34	125,127,137,140	0
2	NAG	B	601	14/15	0.85	0.26	0.33	63,78,95,101	0
3	SO4	E	605	5/5	0.94	0.20	-0.11	59,59,69,75	0
3	SO4	F	603	5/5	0.92	0.22	-0.13	53,66,68,68	0
2	NAG	C	603	14/15	0.71	0.26	-0.34	90,111,122,123	0
3	SO4	C	604	5/5	0.94	0.21	-0.51	38,50,56,61	0
3	SO4	A	603	5/5	0.96	0.19	-0.58	51,52,68,73	0
2	NAG	D	605	14/15	0.79	0.20	-0.72	102,122,135,137	0
2	NAG	C	601	14/15	0.84	0.29	-	83,109,134,154	0
2	NAG	E	601	14/15	0.78	0.35	-	100,128,139,142	0
2	NAG	F	601	14/15	0.65	0.29	-	116,132,146,150	0
2	NAG	C	602	14/15	0.57	0.34	-	156,169,176,185	0
2	NAG	D	604	14/15	0.90	0.29	-	74,90,109,124	0
2	NAG	E	604	14/15	0.77	0.26	-	110,132,162,162	0
2	NAG	B	602	14/15	0.85	0.29	-	79,86,99,100	0
2	NAG	D	601	14/15	0.72	0.22	-	103,132,141,146	0
2	NAG	A	601	14/15	0.73	0.25	-	113,129,140,146	0
2	NAG	D	602	14/15	0.66	0.21	-	130,150,165,168	0
2	NAG	E	602	14/15	0.67	0.29	-	122,149,159,160	0

6.5 Other polymers ⓘ

There are no such residues in this entry.