



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 01:47 PM BST

PDB ID : 5XL5  
Title : The structure of hemagglutinin Q226L mutant from an avian-origin H4N6 influenza virus  
Authors : Song, H.; Qi, J.; Gao, G.F.  
Deposited on : 2017-05-10  
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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.

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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 : 1.8.5 (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

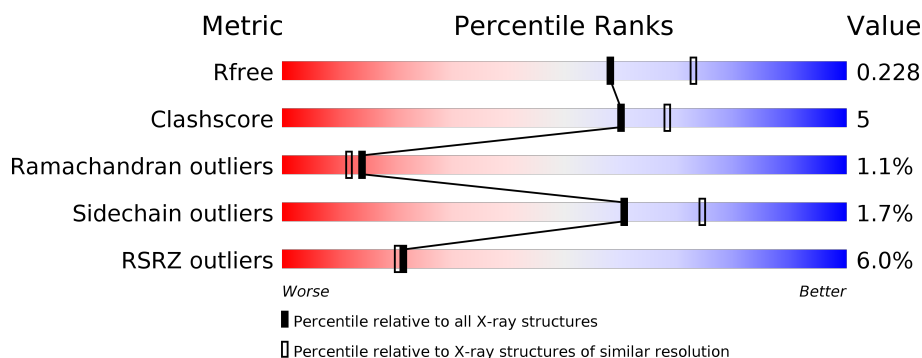
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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))
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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  $\geq 3$ , 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	327	<div> <div>13%</div> <div>81%</div> <div>15%</div> <div>..</div> </div>
1	B	327	<div> <div>3%</div> <div>90%</div> <div>6%</div> <div>..</div> </div>
2	C	176	<div> <div>2%</div> <div>93%</div> <div>..</div> <div>..</div> </div>
2	D	176	<div> <div>2%</div> <div>89%</div> <div>6%</div> <div>..</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8270 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 Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	0	0	0
			2458	1541	436	469	12			
1	B	319	Total	C	N	O	S	0	0	0
			2458	1541	436	469	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	223	LEU	GLN	engineered mutation	UNP A3KF09
B	223	LEU	GLN	engineered mutation	UNP A3KF09

- Molecule 2 is a protein called Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	172	Total	C	N	O	S	0	0	0
			1404	871	250	279	4			
2	D	172	Total	C	N	O	S	0	0	0
			1404	871	250	279	4			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



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

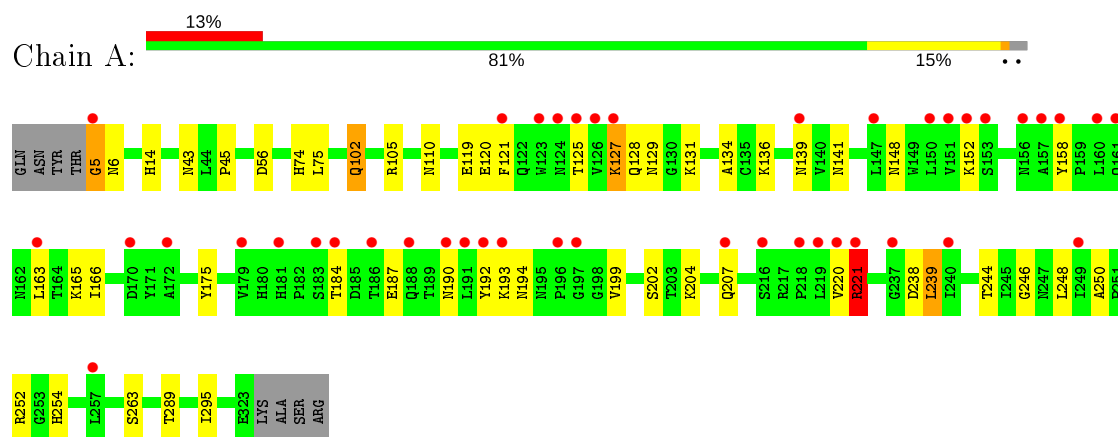
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	63	Total	O	0	0
			63	63		
4	C	126	Total	O	0	0
			126	126		
4	B	156	Total	O	0	0
			156	156		
4	D	145	Total	O	0	0
			145	145		

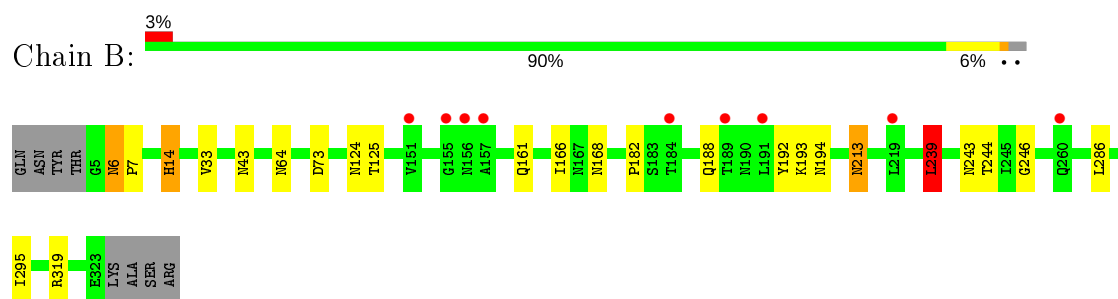
### 3 Residue-property plots [i](#)

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.

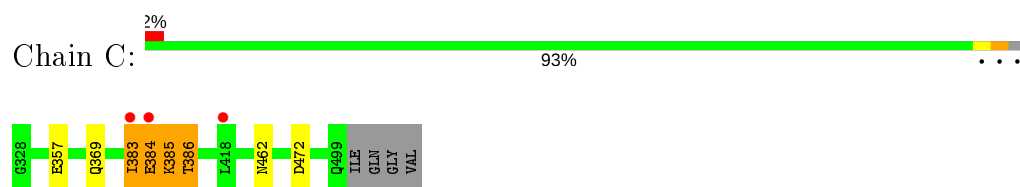
#### • Molecule 1: Hemagglutinin



#### • Molecule 1: Hemagglutinin

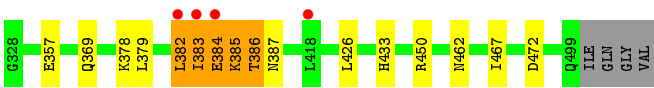


#### • Molecule 2: Hemagglutinin



#### • Molecule 2: Hemagglutinin





## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	100.55Å 100.55Å 685.74Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	33.82 – 2.20 33.82 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (33.82-2.20) 91.3 (33.82-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.62 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.201 , 0.227 0.203 , 0.228	Depositor DCC
$R_{free}$ test set	3500 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.4	Xtriage
Anisotropy	0.457	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 37.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8270	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	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 49.05 % 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 7.8824e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: 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	0.37	0/2512	0.58	1/3420 (0.0%)
1	B	0.42	0/2512	0.59	1/3420 (0.0%)
2	C	0.44	0/1428	0.57	0/1922
2	D	0.47	0/1428	0.61	1/1922 (0.1%)
All	All	0.42	0/7880	0.59	3/10684 (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
2	D	0	2
All	All	0	3

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	239	LEU	CA-CB-CG	6.39	130.00	115.30
1	A	221	ARG	NE-CZ-NH1	-5.86	117.37	120.30
2	D	385	LYS	N-CA-C	5.61	126.15	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	5	GLY	Peptide

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Mol	Chain	Res	Type	Group
2	D	384	GLU	Peptide
2	D	385	LYS	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	2458	0	2412	32	0
1	B	2458	0	2412	19	0
2	C	1404	0	1320	12	0
2	D	1404	0	1320	10	0
3	A	28	0	26	0	0
3	B	28	0	26	0	0
4	A	63	0	0	3	0
4	B	156	0	0	3	0
4	C	126	0	0	2	0
4	D	145	0	0	4	0
All	All	8270	0	7516	71	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 5.

All (71) 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:C:385:LYS:O	2:C:386:THR:CG2	1.79	1.29
2:C:385:LYS:O	2:C:386:THR:HG23	1.01	1.17
1:B:244:THR:HG22	1:B:246:GLY:H	1.29	0.95
2:D:369:GLN:OE1	4:D:701:HOH:O	1.89	0.89
2:C:384:GLU:OE2	2:C:386:THR:HG22	1.73	0.87
2:C:385:LYS:C	2:C:386:THR:HG23	1.94	0.86
1:A:134:ALA:O	1:A:221:ARG:NH1	2.09	0.86
2:D:450:ARG:NH2	4:D:704:HOH:O	2.13	0.81
1:A:102:GLN:HE22	1:A:105:ARG:HH11	1.29	0.79
1:A:244:THR:HG22	1:A:246:GLY:H	1.45	0.79
1:A:56:ASP:OD2	4:A:701:HOH:O	2.00	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:386:THR:O	4:D:702:HOH:O	2.01	0.78
2:D:382:LEU:HD11	2:D:426:LEU:HD21	1.69	0.74
2:D:386:THR:OG1	2:D:387:ASN:N	2.23	0.71
2:D:379:LEU:O	2:D:383:ILE:HG12	1.92	0.70
1:B:73:ASP:OD2	4:B:701:HOH:O	2.10	0.69
2:C:369:GLN:OE1	4:C:601:HOH:O	2.12	0.67
1:A:43:ASN:HB3	1:A:295:ILE:HD13	1.78	0.66
1:B:319:ARG:HG2	1:B:319:ARG:HH11	1.63	0.64
1:B:166:ILE:HA	1:B:239:LEU:HB3	1.81	0.62
1:B:43:ASN:HB3	1:B:295:ILE:HD13	1.84	0.60
1:A:102:GLN:HE22	1:A:105:ARG:NH1	1.97	0.59
1:B:213:ASN:HD22	1:B:213:ASN:N	2.01	0.59
1:A:129:ASN:O	1:A:131:LYS:NZ	2.32	0.58
2:D:357:GLU:OE2	2:D:472:ASP:HB2	2.04	0.58
2:C:385:LYS:C	2:C:386:THR:CG2	2.63	0.58
1:A:148:ASN:HB2	1:A:252:ARG:NE	2.18	0.58
2:C:385:LYS:O	2:C:386:THR:HG22	1.96	0.56
1:B:125:THR:HG22	1:B:125:THR:O	2.05	0.55
1:A:175:TYR:CE2	1:A:254:HIS:HB3	2.43	0.54
2:D:378:LYS:HZ1	2:D:433:HIS:HB3	1.73	0.53
1:B:213:ASN:N	1:B:213:ASN:ND2	2.57	0.52
1:A:192:TYR:O	1:A:194:ASN:N	2.43	0.51
1:B:168:ASN:OD1	4:B:702:HOH:O	2.19	0.51
1:B:73:ASP:CG	4:B:701:HOH:O	2.46	0.50
1:A:120:GLU:HA	1:A:120:GLU:OE1	2.11	0.49
1:B:124:ASN:O	1:B:125:THR:HB	2.12	0.48
1:A:45:PRO:O	4:A:702:HOH:O	2.20	0.48
2:C:357:GLU:OE2	2:C:472:ASP:HB2	2.13	0.48
1:B:182:PRO:HG2	1:B:188:GLN:OE1	2.14	0.47
1:A:121:PHE:HD2	1:A:163:LEU:HD11	1.79	0.47
1:A:102:GLN:NE2	1:A:105:ARG:HD3	2.29	0.47
1:B:161:GLN:O	1:B:243:ASN:HA	2.16	0.46
1:B:213:ASN:HD22	1:B:213:ASN:H	1.63	0.45
1:B:6:ASN:H	1:B:7:PRO:HD3	1.81	0.45
1:A:136:LYS:HE2	1:A:139:ASN:O	2.16	0.45
1:A:204:LYS:HE2	1:A:238:ASP:HA	1.98	0.45
1:A:127:LYS:HD2	1:A:128:GLN:O	2.16	0.45
1:A:110:ASN:HA	1:A:263:SER:O	2.18	0.44
1:A:136:LYS:HE3	1:A:141:ASN:OD1	2.17	0.44
1:B:192:TYR:O	1:B:194:ASN:N	2.50	0.44
2:C:383:ILE:O	2:C:384:GLU:HB2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5:GLY:HA2	4:A:728:HOH:O	2.19	0.43
1:A:184:THR:N	1:A:187:GLU:OE2	2.51	0.43
1:A:244:THR:HG22	1:A:246:GLY:N	2.23	0.43
1:A:202:SER:HA	1:A:207:GLN:HA	2.01	0.42
1:B:7:PRO:HD2	2:D:467:ILE:O	2.20	0.42
1:A:119:GLU:OE2	1:A:165:LYS:HE2	2.20	0.42
1:A:166:ILE:HA	1:A:239:LEU:HB3	2.02	0.42
1:A:289:THR:O	2:C:383:ILE:HG23	2.19	0.42
1:B:6:ASN:H	1:B:7:PRO:CD	2.32	0.42
1:A:199:VAL:HG11	1:A:248:LEU:HD13	2.02	0.41
1:A:158:TYR:O	1:A:194:ASN:ND2	2.53	0.41
2:D:462:ASN:HB2	4:D:772:HOH:O	2.20	0.41
1:A:74:HIS:CE1	1:A:75:LEU:HG	2.56	0.41
1:A:199:VAL:HG22	1:A:244:THR:HG23	2.03	0.41
1:A:148:ASN:N	1:A:250:ALA:O	2.50	0.41
1:A:220:VAL:O	1:A:221:ARG:HG2	2.21	0.40
1:B:14:HIS:CE1	1:B:33:VAL:HG21	2.55	0.40
2:C:462:ASN:HB2	4:C:682:HOH:O	2.21	0.40
2:C:384:GLU:OE2	2:C:386:THR:CG2	2.57	0.40

There are no symmetry-related clashes.

## 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	317/327 (97%)	297 (94%)	16 (5%)	4 (1%)	12	9
1	B	317/327 (97%)	302 (95%)	13 (4%)	2 (1%)	25	26
2	C	170/176 (97%)	159 (94%)	8 (5%)	3 (2%)	8	5
2	D	170/176 (97%)	160 (94%)	8 (5%)	2 (1%)	13	10
All	All	974/1006 (97%)	918 (94%)	45 (5%)	11 (1%)	14	12

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	125	THR
1	A	152	LYS
2	C	384	GLU
2	D	383	ILE
1	A	193	LYS
1	B	6	ASN
2	C	383	ILE
2	D	386	THR
1	B	193	LYS
2	C	386	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	A	275/282 (98%)	269 (98%)	6 (2%)	52	65
1	B	275/282 (98%)	270 (98%)	5 (2%)	59	72
2	C	147/150 (98%)	146 (99%)	1 (1%)	84	91
2	D	147/150 (98%)	145 (99%)	2 (1%)	67	80
All	All	844/864 (98%)	830 (98%)	14 (2%)	60	74

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	HIS
1	A	102	GLN
1	A	127	LYS
1	A	190	ASN
1	A	221	ARG
1	A	239	LEU
2	C	385	LYS
1	B	14	HIS
1	B	64	ASN

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Mol	Chain	Res	Type
1	B	213	ASN
1	B	239	LEU
1	B	286	LEU
2	D	382	LEU
2	D	384	GLU

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	102	GLN
1	B	14	HIS
1	B	64	ASN
1	B	213	ASN

### 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](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

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
3	NAG	B	602	1	14,14,15	0.80	1 (7%)	17,19,21	0.83	1 (5%)
3	NAG	B	601	1	14,14,15	0.21	0	17,19,21	0.64	0
3	NAG	A	602	1	14,14,15	0.31	0	17,19,21	0.59	0
3	NAG	A	601	1	14,14,15	0.52	0	17,19,21	0.52	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	B	602	1	-	2/6/23/26	0/1/1/1
3	NAG	B	601	1	-	2/6/23/26	0/1/1/1
3	NAG	A	602	1	-	0/6/23/26	0/1/1/1
3	NAG	A	601	1	-	1/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	602	NAG	O5-C1	2.64	1.47	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	602	NAG	C1-O5-C5	2.84	116.03	112.19

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	601	NAG	O5-C5-C6-O6
3	B	601	NAG	C4-C5-C6-O6
3	B	602	NAG	O5-C5-C6-O6
3	A	601	NAG	O5-C5-C6-O6
3	B	602	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

## 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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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	319/327 (97%)	0.67	43 (13%) 3 2	27, 72, 120, 139	0
1	B	319/327 (97%)	-0.22	9 (2%) 53 51	20, 41, 66, 100	0
2	C	172/176 (97%)	0.09	3 (1%) 70 68	26, 35, 64, 112	0
2	D	172/176 (97%)	-0.07	4 (2%) 60 58	21, 30, 46, 119	0
All	All	982/1006 (97%)	0.15	59 (6%) 21 20	20, 41, 99, 139	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	383	ILE	10.2
1	A	123	TRP	7.1
1	A	157	ALA	5.9
1	A	197	GLY	5.8
1	A	192	TYR	5.4
1	A	151	VAL	5.1
1	A	184	THR	4.8
2	C	383	ILE	4.2
1	A	158	TYR	4.2
2	D	384	GLU	4.0
1	A	191	LEU	4.0
1	A	188	GLN	3.9
1	A	156	ASN	3.5
1	B	155	GLY	3.5
1	A	181	HIS	3.2
1	A	207	GLN	3.0
1	A	196	PRO	3.0
1	A	160	LEU	2.9
1	A	5	GLY	2.9
1	A	183	SER	2.9
1	A	218	PRO	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	127	LYS	2.9
1	A	126	VAL	2.8
1	A	125	THR	2.8
1	A	139	ASN	2.7
2	C	418	LEU	2.7
1	B	151	VAL	2.7
1	A	170	ASP	2.6
1	A	147	LEU	2.6
1	A	240	ILE	2.6
1	A	249	ILE	2.6
1	B	157	ALA	2.5
1	A	221	ARG	2.5
1	A	186	THR	2.5
2	D	418	LEU	2.5
1	A	163	LEU	2.5
1	A	153	SER	2.5
1	A	219	LEU	2.5
1	A	172	ALA	2.4
1	A	220	VAL	2.4
1	B	189	THR	2.4
2	D	382	LEU	2.4
1	A	124	ASN	2.3
1	A	152	LYS	2.3
1	B	184	THR	2.3
2	C	384	GLU	2.3
1	A	121	PHE	2.3
1	A	179	VAL	2.2
1	A	193	LYS	2.1
1	A	161	GLN	2.1
1	A	150	LEU	2.1
1	B	191	LEU	2.1
1	A	190	ASN	2.1
1	A	257	LEU	2.0
1	A	237	GLY	2.0
1	B	156	ASN	2.0
1	B	219	LEU	2.0
1	B	260	GLN	2.0
1	A	216	SER	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 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
3	NAG	A	601	14/15	0.81	0.33	108,116,122,125	0
3	NAG	B	601	14/15	0.86	0.27	92,95,98,99	0
3	NAG	B	602	14/15	0.91	0.12	28,33,39,40	0
3	NAG	A	602	14/15	0.93	0.16	40,48,51,54	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.