



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 25, 2025 – 06:08 PM JST

PDB ID : 9JEB / pdb_00009jeb
Title : Crystal structure of SARS-CoV-2 RBD in complex with a neutralizing anti-body scFv N1
Authors : Guo, K.; Zhang, N.; Guo, Y.
Deposited on : 2024-09-03
Resolution : 2.30 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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0rc1
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.45.1

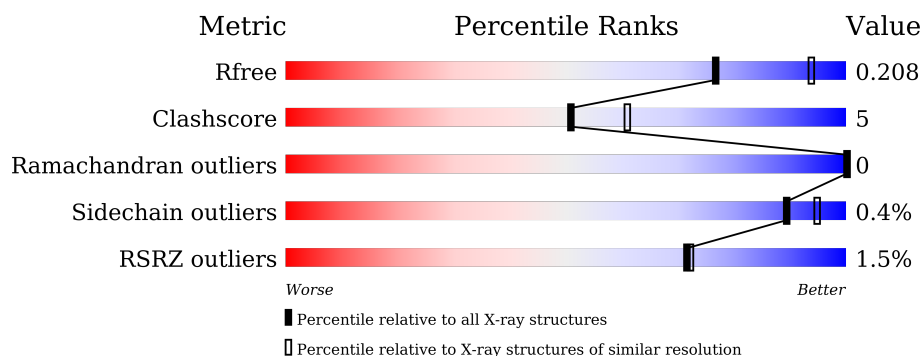
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	5963 (2.30-2.30)
Clashscore	180529	6698 (2.30-2.30)
Ramachandran outliers	177936	6640 (2.30-2.30)
Sidechain outliers	177891	6640 (2.30-2.30)
RSRZ outliers	164620	5963 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	247	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, orange 1%, orange 80%, yellow 80%, yellow 93%, green 93%, green 100%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 80% 13% • 6% </div> </div>
1	C	247	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 0%, green 81%, yellow 81%, yellow 94%, green 94%, green 100%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 81% 13% 6% </div> </div>
2	B	197	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 3%, orange 3%, orange 91%, yellow 91%, yellow 99%, green 99%, green 100%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 3% 91% 8% •• </div> </div>
2	D	197	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 2%, orange 2%, orange 91%, yellow 91%, yellow 99%, green 99%, green 100%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 2% 91% 9% • </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7532 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.

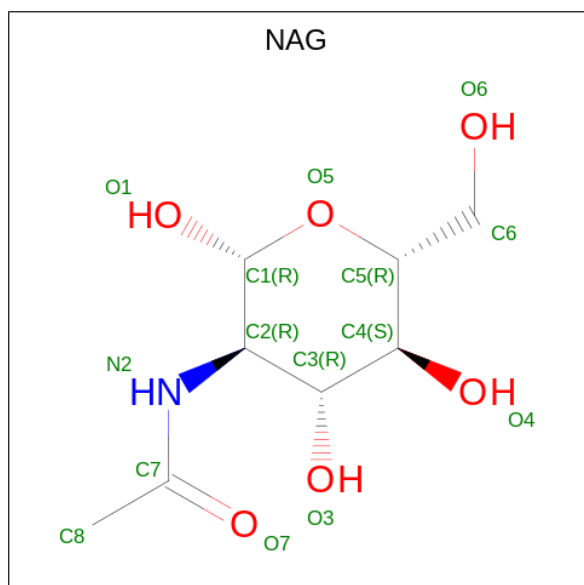
- Molecule 1 is a protein called N1 scFv.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	232	Total	C	N	O	S	0	0	0
			1790	1132	299	353	6			
1	C	233	Total	C	N	O	S	0	0	0
			1796	1135	300	355	6			

- Molecule 2 is a protein called Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	196	Total	C	N	O	S	0	0	0
			1552	995	259	290	8			
2	D	197	Total	C	N	O	S	0	0	0
			1561	1001	261	291	8			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

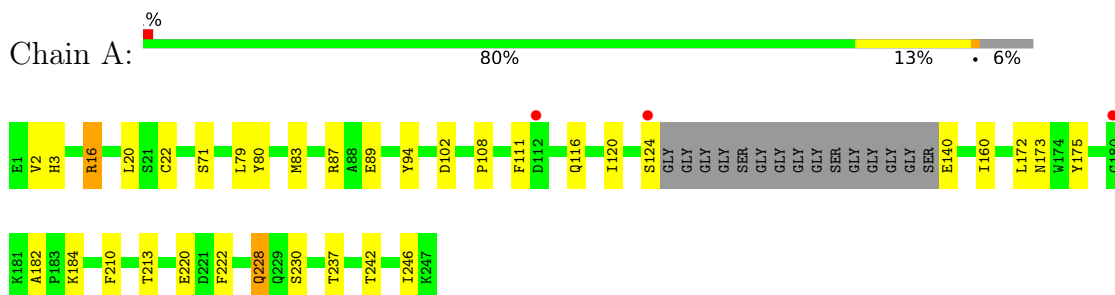
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	246	Total	O	0	0
			246	246		
4	B	175	Total	O	0	0
			175	175		
4	C	210	Total	O	0	0
			210	210		
4	D	188	Total	O	0	0
			188	188		

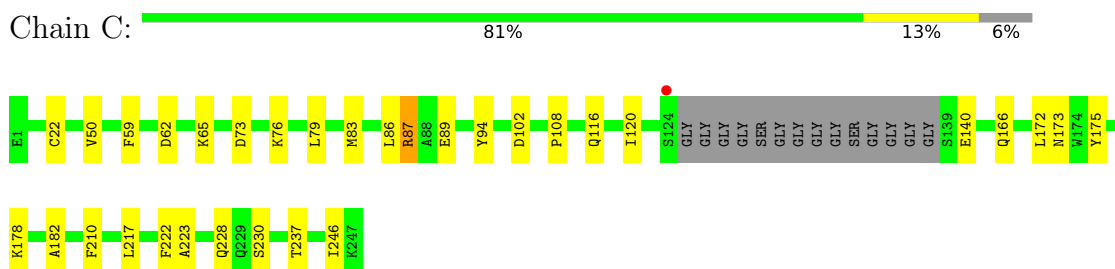
3 Residue-property plots [i](#)

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.

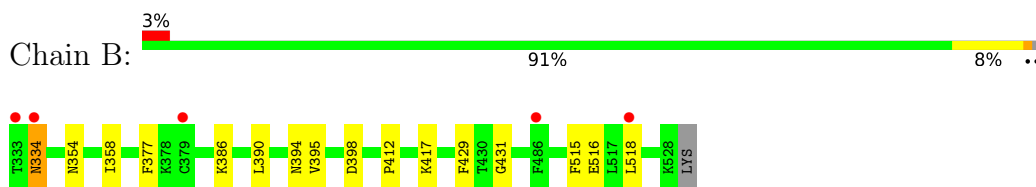
- Molecule 1: N1 scFv



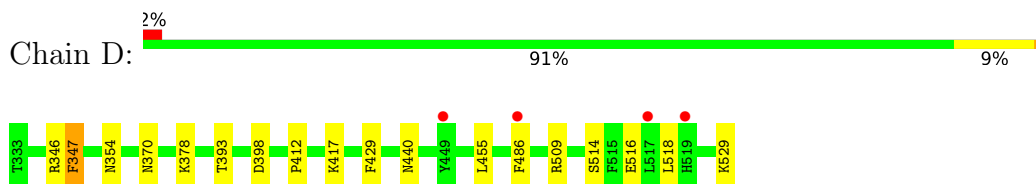
- Molecule 1: N1 scFv



- Molecule 2: Spike protein S1



- Molecule 2: Spike protein S1



4 Data and refinement statistics

Property	Value	Source
Space group	P 62	Depositor
Cell constants a, b, c, α , β , γ	190.40Å 190.40Å 60.14Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	32.98 – 2.30 32.98 – 2.30	Depositor EDS
% Data completeness (in resolution range)	90.5 (32.98-2.30) 90.5 (32.98-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.64 (at 2.29Å)	Xtriage
Refinement program	PHENIX (1.18.2_3874: ???)	Depositor
R, R_{free}	0.178 , 0.206 0.179 , 0.208	Depositor DCC
R_{free} test set	2024 reflections (3.63%)	wwPDB-VP
Wilson B-factor (Å ²)	30.4	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.328 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7532	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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	A	0.51	2/1835 (0.1%)	0.79	4/2489 (0.2%)
1	C	0.44	0/1841	0.74	1/2497 (0.0%)
2	B	0.39	0/1596	0.71	0/2172
2	D	0.45	0/1605	0.69	1/2183 (0.0%)
All	All	0.45	2/6877 (0.0%)	0.73	6/9341 (0.1%)

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	B	0	1
2	D	0	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	111	PHE	C-O	-5.92	1.15	1.23
1	A	2	VAL	C-O	-5.04	1.18	1.24

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	16	ARG	CD-NE-CZ	9.55	137.77	124.40
1	A	16	ARG	NE-CZ-NH1	-7.64	113.86	121.50
1	A	87	ARG	CB-CG-CD	-6.36	96.67	111.30
1	A	228	GLN	CA-CB-CG	5.65	125.41	114.10

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	87	ARG	CG-CD-NE	5.65	124.44	112.00
2	D	486	PHE	CA-C-O	-5.34	114.83	120.92

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	16	ARG	Sidechain
2	B	377	PHE	Sidechain
2	D	347	PHE	Sidechain

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	1790	0	1718	22	0
1	C	1796	0	1723	21	0
2	B	1552	0	1473	12	0
2	D	1561	0	1485	12	0
3	D	14	0	13	0	0
4	A	246	0	0	7	0
4	B	175	0	0	1	1
4	C	210	0	0	3	0
4	D	188	0	0	5	1
All	All	7532	0	6412	67	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 5.

All (67) 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:B:334:ASN:H	2:B:334:ASN:HD22	1.16	0.91
2:D:393:THR:OG1	4:D:701:HOH:O	2.02	0.77
2:D:370:ASN:ND2	4:D:702:HOH:O	2.13	0.73

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:HIS:ND1	4:A:303:HOH:O	2.23	0.72
2:D:378:LYS:HE3	4:D:862:HOH:O	1.90	0.71
2:D:516:GLU:HG2	2:D:518:LEU:HG	1.73	0.70
1:A:175:TYR:HE2	1:A:228:GLN:HG3	1.56	0.69
1:C:166:GLN:HG3	4:C:310:HOH:O	1.94	0.67
1:C:175:TYR:HE2	1:C:228:GLN:HG3	1.60	0.67
2:B:334:ASN:H	2:B:334:ASN:ND2	1.93	0.66
2:B:334:ASN:HD22	2:B:334:ASN:N	1.93	0.65
1:A:213:THR:HG21	4:A:528:HOH:O	1.98	0.62
1:A:71:SER:OG	4:A:301:HOH:O	2.12	0.62
1:A:173:ASN:HD22	1:A:228:GLN:NE2	1.99	0.61
1:A:220:GLU:OE2	4:A:302:HOH:O	2.15	0.60
1:A:213:THR:HG23	4:A:435:HOH:O	2.01	0.59
2:B:412:PRO:HG3	2:B:429:PHE:HB3	1.87	0.56
1:C:94:TYR:CD1	1:C:120:ILE:HD13	2.41	0.56
2:D:417:LYS:HE2	2:D:455:LEU:HD12	1.88	0.55
1:C:175:TYR:HE2	1:C:228:GLN:CG	2.19	0.55
1:C:87:ARG:HG3	1:C:89:GLU:OE1	2.07	0.55
1:A:20:LEU:HG	1:A:83:MET:HE2	1.89	0.55
1:C:173:ASN:HD22	1:C:228:GLN:NE2	2.05	0.54
1:A:228:GLN:HB2	1:A:237:THR:O	2.08	0.54
1:C:73:ASP:OD2	1:C:76:LYS:HE3	2.09	0.52
1:C:102:ASP:HB2	4:C:359:HOH:O	2.09	0.52
1:C:62:ASP:OD1	1:C:65:LYS:NZ	2.43	0.52
2:D:429:PHE:HE1	2:D:514:SER:HB3	1.75	0.51
1:A:184:LYS:HG2	4:A:381:HOH:O	2.11	0.50
1:A:160:ILE:HG12	1:A:242:THR:HG21	1.95	0.49
1:A:116:GLN:HA	1:A:182:ALA:HB2	1.94	0.49
1:A:102:ASP:HB2	4:A:358:HOH:O	2.13	0.49
2:D:412:PRO:HG3	2:D:429:PHE:HB3	1.95	0.48
1:A:71:SER:HB3	1:A:80:TYR:HB2	1.95	0.48
2:B:417:LYS:NZ	4:B:607:HOH:O	2.47	0.47
2:B:334:ASN:ND2	2:B:334:ASN:N	2.58	0.47
1:A:108:PRO:HB3	1:A:230:SER:O	2.15	0.46
2:B:394:ASN:HB2	2:B:516:GLU:OE1	2.15	0.46
1:C:22:CYS:HB3	1:C:79:LEU:HB3	1.96	0.46
1:C:172:LEU:HD22	1:C:210:PHE:CB	2.45	0.46
1:A:222:PHE:HB3	1:A:246:ILE:HG12	1.97	0.46
1:C:83:MET:HE2	1:C:86:LEU:HD21	1.97	0.46
1:A:172:LEU:HD22	1:A:210:PHE:CB	2.45	0.46
2:B:516:GLU:HG2	2:B:518:LEU:HG	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:140:GLU:HG3	1:C:237:THR:HG21	1.97	0.45
1:C:222:PHE:HB3	1:C:246:ILE:HG12	1.98	0.45
1:C:116:GLN:HA	1:C:182:ALA:HB2	1.99	0.45
1:C:228:GLN:HB2	1:C:237:THR:O	2.17	0.44
2:D:347:PHE:CE2	2:D:509:ARG:HB3	2.52	0.44
1:A:94:TYR:CD1	1:A:120:ILE:HD13	2.53	0.44
1:C:108:PRO:HB3	1:C:230:SER:O	2.18	0.44
2:D:440:ASN:ND2	4:D:707:HOH:O	2.45	0.44
1:C:87:ARG:HH21	1:C:87:ARG:HD2	1.73	0.43
2:B:354:ASN:O	2:B:398:ASP:HA	2.19	0.43
2:D:346:ARG:NH2	4:D:711:HOH:O	2.50	0.43
2:B:386:LYS:O	2:B:390:LEU:HG	2.18	0.43
1:A:22:CYS:HB3	1:A:79:LEU:HB3	2.00	0.42
1:A:172:LEU:HD22	1:A:210:PHE:CG	2.55	0.42
2:B:358:ILE:HB	2:B:395:VAL:HB	2.02	0.42
2:B:431:GLY:HA2	2:B:515:PHE:CD2	2.55	0.42
1:A:175:TYR:HE2	1:A:228:GLN:CG	2.29	0.41
1:A:89:GLU:H	1:A:89:GLU:CD	2.28	0.41
1:C:217:LEU:HD21	1:C:246:ILE:CD1	2.51	0.41
2:D:354:ASN:O	2:D:398:ASP:HA	2.21	0.41
4:C:355:HOH:O	2:D:529:LYS:HE2	2.20	0.40
1:C:50:VAL:HG12	1:C:59:PHE:HB2	2.04	0.40
1:C:178:LYS:HG2	1:C:223:ALA:HB2	2.02	0.40

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)
4:B:746:HOH:O	4:D:880:HOH:O[4_456]	1.81	0.39

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	228/247 (92%)	223 (98%)	5 (2%)	0	100	100
1	C	229/247 (93%)	223 (97%)	6 (3%)	0	100	100
2	B	194/197 (98%)	189 (97%)	5 (3%)	0	100	100
2	D	195/197 (99%)	190 (97%)	5 (3%)	0	100	100
All	All	846/888 (95%)	825 (98%)	21 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	197/200 (98%)	195 (99%)	2 (1%)	73	85
1	C	198/200 (99%)	198 (100%)	0	100	100
2	B	169/170 (99%)	168 (99%)	1 (1%)	84	92
2	D	170/170 (100%)	170 (100%)	0	100	100
All	All	734/740 (99%)	731 (100%)	3 (0%)	89	95

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	124	SER
1	A	140	GLU
2	B	334	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	82	GLN
1	A	84	ASN
1	A	176	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	177	GLN
1	A	228	GLN
2	B	334	ASN
2	B	354	ASN
2	B	474	GLN
1	C	39	GLN
1	C	53	HIS
1	C	77	ASN
1	C	177	GLN
1	C	228	GLN
2	D	498	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	D	601	2	14,14,15	1.45	2 (14%)	17,19,21	1.62	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	601	2	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	601	NAG	C2-N2	-2.86	1.41	1.46
3	D	601	NAG	O7-C7	-2.27	1.18	1.23

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	601	NAG	O5-C5-C6	4.51	114.28	107.20

There are no chirality outliers.

There are no torsion outliers.

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 [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	232/247 (93%)	-0.29	3 (1%) 74 75	21, 32, 47, 71	0
1	C	233/247 (94%)	-0.20	1 (0%) 89 89	23, 34, 50, 67	0
2	B	196/197 (99%)	-0.04	5 (2%) 57 58	25, 37, 56, 69	0
2	D	197/197 (100%)	-0.10	4 (2%) 64 66	24, 35, 56, 71	0
All	All	858/888 (96%)	-0.16	13 (1%) 71 72	21, 34, 54, 71	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	486	PHE	3.3
2	B	379	CYS	3.0
1	A	180	GLY	3.0
1	A	124	SER	2.8
2	D	519	HIS	2.8
2	B	486	PHE	2.8
2	B	518	LEU	2.7
1	C	124	SER	2.6
2	D	449	TYR	2.4
2	B	334	ASN	2.2
1	A	112	ASP	2.2
2	D	517	LEU	2.2
2	B	333	THR	2.1

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 oligosaccharides 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, 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
3	NAG	D	601	14/15	0.95	0.07	36,40,47,49	0

6.5 Other polymers [i](#)

There are no such residues in this entry.