



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:35 PM UTC

PDB ID : 5XL2 / pdb_00005xl2
Title : The structure of hemagglutinin from a swine-origin H4N6 influenza virus
Authors : Song, H.; Qi, J.; Gao, G.F.
Deposited on : 2017-05-10
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 Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

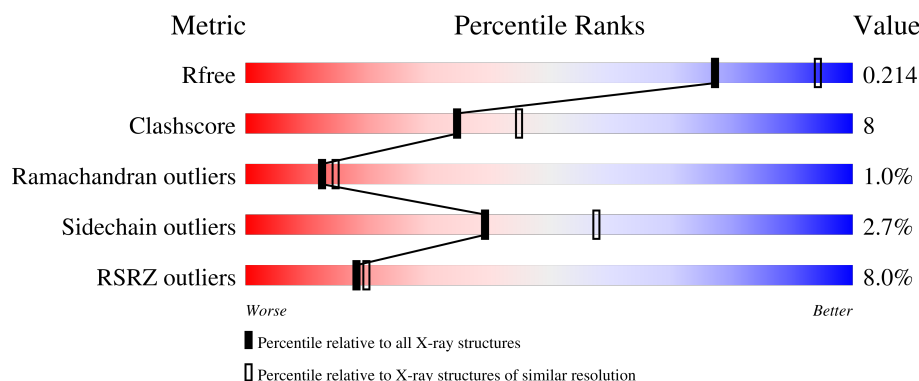
1 Overall quality at a glance

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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (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	503	3% 86% 10% ..
1	B	503	7% 81% 14% ..
1	C	503	14% 78% 14% ...

2 Entry composition [i](#)

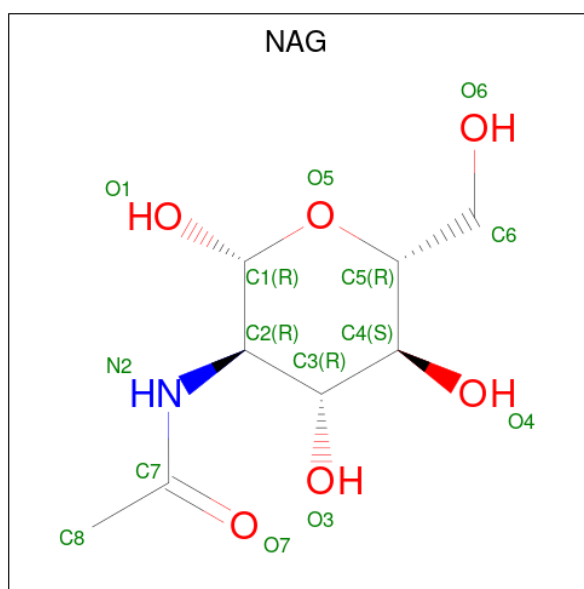
There are 3 unique types of molecules in this entry. The entry contains 12325 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	486	Total	C	N	O	S	0	0	0
			3841	2398	680	748	15			
1	B	486	Total	C	N	O	S	0	0	0
			3843	2400	680	748	15			
1	C	484	Total	C	N	O	S	0	0	0
			3830	2391	678	746	15			

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



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

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
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		

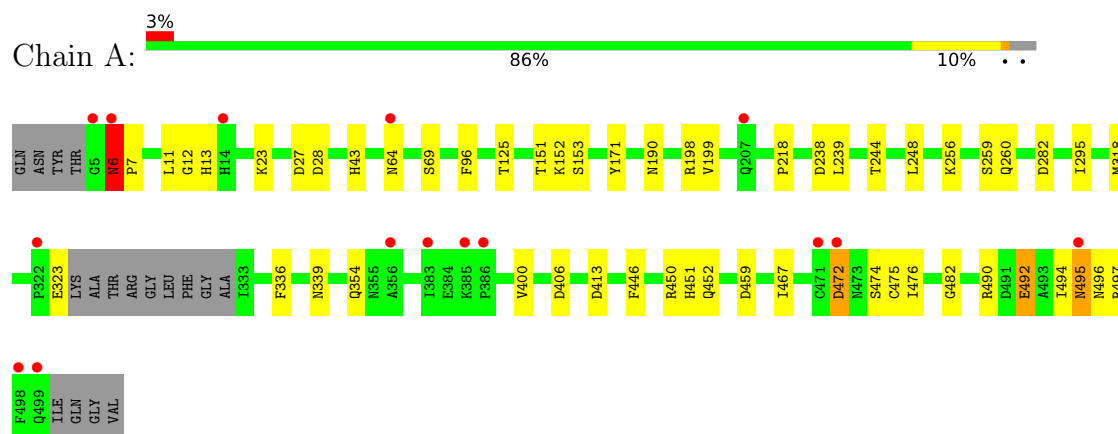
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	274	Total	O	0	0
			274	274		
3	B	240	Total	O	0	0
			240	240		
3	C	213	Total	O	0	0
			213	213		

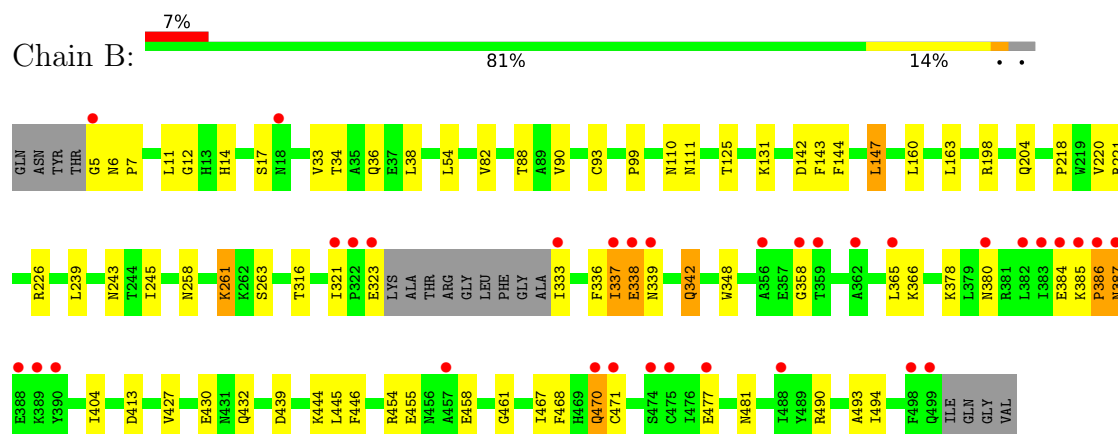
3 Residue-property plots

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.

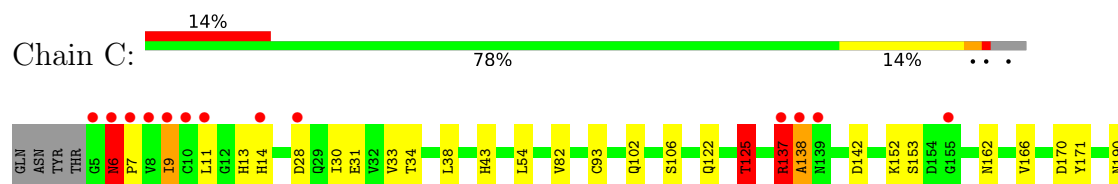
• Molecule 1: Hemagglutinin

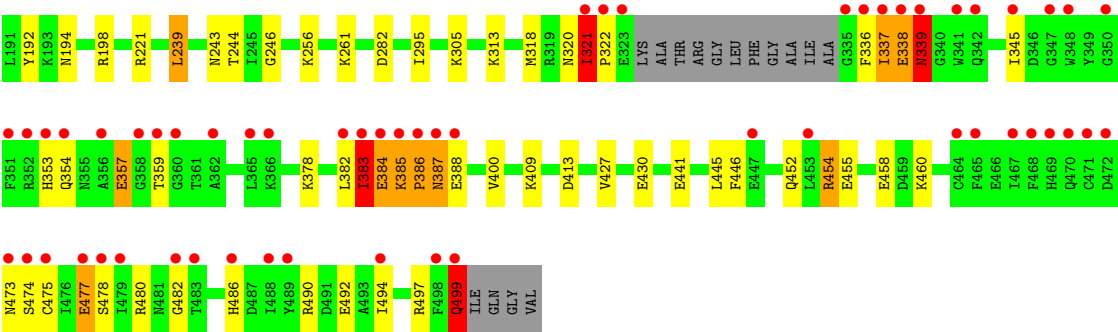


• Molecule 1: Hemagglutinin



• Molecule 1: Hemagglutinin





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.42Å 120.56Å 132.28Å 90.00° 92.29° 90.00°	Depositor
Resolution (Å)	35.57 – 2.30 35.57 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.5 (35.57-2.30) 99.7 (35.57-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.09 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.184 , 0.213 0.187 , 0.214	Depositor DCC
R_{free} test set	4563 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	1.122	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.023 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12325	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% 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

5.1 Standard geometry

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	0.40	0/3923	0.80	2/5323 (0.0%)
1	B	0.43	0/3925	0.79	6/5327 (0.1%)
1	C	0.41	0/3912	0.81	12/5309 (0.2%)
All	All	0.41	0/11760	0.80	20/15959 (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	C	0	1

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	6	ASN	CA-C-N	-9.72	109.49	119.99
1	A	6	ASN	C-N-CA	-9.72	109.49	119.99
1	C	475	CYS	N-CA-C	-7.70	104.02	113.41
1	C	6	ASN	CA-C-N	-7.25	112.51	119.76
1	C	6	ASN	C-N-CA	-7.25	112.51	119.76
1	C	339	ASN	N-CA-C	-6.42	97.12	110.80
1	C	385	LYS	N-CA-C	6.37	123.90	109.81
1	B	358	GLY	N-CA-C	6.24	118.44	110.20
1	B	321	ILE	CA-C-N	6.09	126.03	119.76
1	B	321	ILE	C-N-CA	6.09	126.03	119.76
1	C	125	THR	CB-CA-C	-5.86	98.77	110.42
1	C	499	GLN	CA-CB-CG	5.65	125.40	114.10
1	C	454	ARG	CB-CA-C	-5.61	110.09	116.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	321	ILE	CA-C-N	5.58	125.53	119.78
1	C	321	ILE	C-N-CA	5.58	125.53	119.78
1	B	454	ARG	CB-CA-C	-5.54	110.20	116.63
1	C	339	ASN	CA-C-N	-5.54	116.71	122.69
1	C	339	ASN	C-N-CA	-5.54	116.71	122.69
1	B	342	GLN	N-CA-C	5.13	119.00	112.34
1	B	338	GLU	N-CA-C	5.07	116.39	110.41

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	384	GLU	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	3841	0	3690	47	0
1	B	3843	0	3698	54	0
1	C	3830	0	3682	75	0
2	A	28	0	26	0	0
2	B	28	0	26	0	0
2	C	28	0	26	2	0
3	A	274	0	0	6	0
3	B	240	0	0	12	0
3	C	213	0	0	12	0
All	All	12325	0	11148	171	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 8.

All (171) 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:C:137:ARG:NE	1:C:142:ASP:OD2	1.57	1.36

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:499:GLN:HE21	1:C:499:GLN:HA	1.21	1.02
1:A:495:ASN:C	1:A:495:ASN:HD22	1.76	0.93
1:B:439:ASP:OD1	3:B:702:HOH:O	1.89	0.91
1:A:244:THR:HG21	1:A:248:LEU:HB2	1.53	0.90
1:C:122:GLN:OE1	3:C:701:HOH:O	1.90	0.89
1:B:413:ASP:OD1	3:B:703:HOH:O	1.89	0.89
1:A:354:GLN:OE1	3:A:701:HOH:O	1.95	0.85
1:B:385:LYS:HB3	1:B:386:PRO:HD2	1.59	0.85
1:B:384:GLU:HG3	1:B:385:LYS:H	1.42	0.84
1:C:244:THR:HG22	1:C:246:GLY:H	1.45	0.82
1:A:495:ASN:O	1:A:495:ASN:ND2	2.12	0.81
1:B:14:HIS:HD2	1:B:348:TRP:HA	1.44	0.79
1:C:198:ARG:NH2	1:C:243:ASN:OD1	2.14	0.79
1:C:137:ARG:CD	1:C:142:ASP:OD2	2.31	0.78
1:C:492:GLU:OE1	3:C:702:HOH:O	1.99	0.78
1:A:413:ASP:OD1	3:A:702:HOH:O	2.02	0.76
1:A:406:ASP:OD1	3:A:703:HOH:O	2.03	0.76
1:B:198:ARG:NH2	1:B:243:ASN:OD1	2.19	0.74
1:B:455:GLU:OE2	3:B:705:HOH:O	2.05	0.74
1:A:451:HIS:ND1	1:B:461:GLY:HA2	2.04	0.73
1:C:441:GLU:O	3:C:704:HOH:O	2.06	0.72
2:C:602:NAG:O7	3:C:705:HOH:O	2.08	0.72
1:C:474:SER:HA	1:C:477:GLU:HG2	1.71	0.71
1:C:93:CYS:O	1:C:221:ARG:NH1	2.24	0.71
1:B:14:HIS:CD2	1:B:348:TRP:HA	2.25	0.71
1:C:413:ASP:OD1	3:C:706:HOH:O	2.09	0.71
1:C:137:ARG:HE	1:C:142:ASP:CG	1.94	0.70
1:C:490:ARG:O	1:C:494:ILE:HG12	1.92	0.69
1:C:499:GLN:HA	1:C:499:GLN:NE2	2.04	0.68
1:B:12:GLY:HA2	1:B:336:PHE:HB3	1.74	0.68
1:A:495:ASN:C	1:A:495:ASN:ND2	2.51	0.68
1:B:481:ASN:ND2	3:B:704:HOH:O	2.01	0.67
1:C:11:LEU:HD13	1:C:446:PHE:HA	1.76	0.67
1:C:386:PRO:O	1:C:387:ASN:O	2.12	0.67
1:C:354:GLN:HB2	1:C:359:THR:HG22	1.75	0.66
1:B:11:LEU:HD13	1:B:446:PHE:HA	1.78	0.66
1:A:492:GLU:O	1:A:496:ASN:ND2	2.28	0.64
1:C:385:LYS:NZ	3:C:703:HOH:O	2.05	0.64
1:C:388:GLU:N	1:C:388:GLU:OE1	2.30	0.64
1:C:9:ILE:HD12	1:C:353:HIS:HB3	1.78	0.64
1:B:218:PRO:O	1:B:226:ARG:NH2	2.29	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:378:LYS:NZ	1:B:430:GLU:O	2.31	0.64
1:A:199:VAL:HA	1:A:244:THR:HG22	1.81	0.63
1:A:472:ASP:HB3	1:A:474:SER:H	1.64	0.63
1:C:338:GLU:O	1:C:339:ASN:ND2	2.32	0.62
1:B:384:GLU:CG	1:B:385:LYS:H	2.13	0.62
1:A:497:ARG:NH2	1:C:455:GLU:OE2	2.32	0.62
1:C:452:GLN:NE2	1:C:482:GLY:HA2	2.14	0.62
1:A:450:ARG:NH2	1:A:459:ASP:OD2	2.33	0.61
1:B:490:ARG:NH2	1:C:458:GLU:OE2	2.34	0.60
1:C:125:THR:O	1:C:125:THR:HG22	1.99	0.60
1:C:33:VAL:HG12	1:C:34:THR:HG23	1.83	0.59
1:B:220:VAL:HG23	1:B:226:ARG:NH2	2.17	0.59
1:B:323:GLU:HG3	1:B:339:ASN:HD22	1.67	0.59
1:A:452:GLN:NE2	1:A:482:GLY:HA2	2.18	0.58
1:A:64:ASN:ND2	1:A:96:PHE:CZ	2.72	0.58
1:B:384:GLU:HG3	1:B:385:LYS:N	2.17	0.58
1:A:450:ARG:HG2	1:A:451:HIS:HD2	1.69	0.57
1:C:171:TYR:CZ	1:C:256:LYS:HD2	2.39	0.57
1:C:321:ILE:HG22	1:C:322:PRO:HD2	1.86	0.57
1:B:432:GLN:NE2	3:B:701:HOH:O	1.88	0.57
1:A:64:ASN:HB3	1:A:69:SER:OG	2.05	0.57
1:C:357:GLU:OE2	1:C:473:ASN:N	2.26	0.57
1:B:261:LYS:HD2	1:C:400:VAL:HG11	1.87	0.55
1:B:221:ARG:NH1	3:B:708:HOH:O	2.19	0.55
1:B:131:LYS:HD2	1:B:142:ASP:HA	1.88	0.55
1:C:102:GLN:CD	1:C:102:GLN:H	2.14	0.55
1:C:322:PRO:O	1:C:339:ASN:HB3	2.07	0.55
1:A:11:LEU:HD13	1:A:446:PHE:HA	1.89	0.54
1:B:99:PRO:O	3:B:707:HOH:O	2.19	0.54
1:C:338:GLU:HG2	1:C:339:ASN:HB2	1.89	0.54
1:A:64:ASN:ND2	1:A:96:PHE:HZ	2.06	0.54
1:A:218:PRO:HG3	1:C:239:LEU:HD22	1.90	0.54
1:C:43:HIS:HB3	1:C:295:ILE:HD13	1.91	0.53
1:C:378:LYS:NZ	1:C:430:GLU:O	2.43	0.52
1:A:323:GLU:HA	1:A:339:ASN:HD22	1.74	0.52
1:B:198:ARG:HG3	1:B:245:ILE:HG12	1.91	0.52
1:A:400:VAL:HG11	1:C:261:LYS:HD2	1.91	0.51
1:A:171:TYR:CZ	1:A:256:LYS:HD3	2.45	0.51
1:C:474:SER:HA	1:C:477:GLU:CG	2.39	0.51
1:C:162:ASN:HD22	2:C:601:NAG:H83	1.75	0.50
1:B:337:ILE:HG22	1:B:338:GLU:HG3	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:31:GLU:HG2	1:C:320:ASN:HB3	1.92	0.50
1:C:28:ASP:OD1	1:C:28:ASP:N	2.35	0.50
1:A:125:THR:HG22	1:A:125:THR:O	2.11	0.49
1:B:125:THR:HG22	1:B:125:THR:O	2.11	0.49
1:C:54:LEU:HD22	1:C:82:VAL:HB	1.94	0.49
1:C:409:LYS:NZ	3:C:726:HOH:O	2.45	0.49
1:A:6:ASN:H	1:A:7:PRO:HD2	1.77	0.49
1:A:472:ASP:HB2	1:A:475:CYS:H	1.78	0.49
1:C:6:ASN:H	1:C:7:PRO:CD	2.25	0.49
1:C:305:LYS:NZ	3:C:715:HOH:O	2.31	0.49
1:B:258:ASN:O	1:B:261:LYS:HG2	2.13	0.48
1:B:36:GLN:NE2	3:B:727:HOH:O	2.45	0.48
1:A:7:PRO:HG2	1:A:467:ILE:HB	1.95	0.48
1:A:43:HIS:HB3	1:A:295:ILE:HD13	1.96	0.48
1:B:88:THR:OG1	3:B:709:HOH:O	2.20	0.47
1:C:455:GLU:O	1:C:497:ARG:NH1	2.40	0.47
1:A:27:ASP:OD2	1:A:27:ASP:N	2.47	0.47
1:A:238:ASP:OD1	1:A:239:LEU:N	2.42	0.47
1:B:458:GLU:HB2	1:B:468:PHE:HE2	1.79	0.47
1:A:198:ARG:O	1:A:244:THR:HG22	2.14	0.47
1:B:142:ASP:CG	1:B:143:PHE:H	2.22	0.47
1:B:468:PHE:O	1:B:493:ALA:HA	2.14	0.47
1:B:490:ARG:O	1:B:494:ILE:HG12	2.15	0.47
1:C:386:PRO:O	1:C:387:ASN:C	2.58	0.47
1:A:6:ASN:H	1:A:7:PRO:CD	2.28	0.47
1:B:93:CYS:O	1:B:221:ARG:NH1	2.46	0.47
1:C:388:GLU:H	1:C:388:GLU:CD	2.23	0.47
1:A:23:LYS:NZ	1:A:28:ASP:HB2	2.30	0.46
1:C:38:LEU:HD12	1:C:427:VAL:HG22	1.95	0.46
1:B:5:GLY:HA2	1:B:468:PHE:HA	1.97	0.46
1:C:137:ARG:HB3	1:C:138:ALA:H	1.61	0.46
1:C:452:GLN:HE22	1:C:482:GLY:HA2	1.79	0.46
1:B:110:ASN:HA	1:B:263:SER:O	2.15	0.46
1:B:7:PRO:HD2	1:B:467:ILE:O	2.16	0.46
1:C:385:LYS:HG3	1:C:385:LYS:O	2.16	0.46
1:C:11:LEU:HD22	1:C:445:LEU:HG	1.97	0.45
1:A:23:LYS:HE2	3:A:834:HOH:O	2.17	0.45
1:A:23:LYS:HZ2	1:A:28:ASP:HB2	1.82	0.45
1:C:106:SER:OG	3:C:709:HOH:O	2.21	0.44
1:C:337:ILE:O	1:C:338:GLU:C	2.60	0.44
1:A:490:ARG:O	1:A:494:ILE:HG12	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:54:LEU:HD22	1:B:82:VAL:HB	1.98	0.44
1:C:383:ILE:HD12	1:C:383:ILE:O	2.17	0.44
1:C:477:GLU:HG3	1:C:478:SER:H	1.83	0.44
1:C:313:LYS:O	3:C:708:HOH:O	2.21	0.44
1:B:380:ASN:ND2	3:B:715:HOH:O	2.50	0.43
1:C:477:GLU:HG3	1:C:478:SER:N	2.32	0.43
1:B:142:ASP:OD1	1:B:143:PHE:N	2.51	0.43
1:C:137:ARG:HD2	1:C:142:ASP:OD2	2.13	0.43
1:C:460:LYS:HE2	1:C:460:LYS:HB3	1.85	0.43
1:A:6:ASN:ND2	3:A:704:HOH:O	2.09	0.43
1:C:192:TYR:O	1:C:194:ASN:N	2.47	0.43
1:C:477:GLU:O	1:C:480:ARG:N	2.50	0.43
1:A:152:LYS:HE2	1:A:190:ASN:O	2.19	0.43
1:B:337:ILE:O	1:B:338:GLU:HG2	2.19	0.43
1:C:13:HIS:HB2	1:C:318:MET:SD	2.59	0.43
1:C:244:THR:HG22	1:C:246:GLY:N	2.23	0.43
1:A:12:GLY:HA2	1:A:336:PHE:HB3	2.01	0.43
1:A:323:GLU:OE1	3:A:706:HOH:O	2.21	0.43
1:A:452:GLN:HE21	1:A:482:GLY:HA2	1.84	0.43
1:C:152:LYS:HE2	1:C:190:ASN:O	2.18	0.43
1:C:166:VAL:HA	1:C:239:LEU:HB3	2.01	0.43
1:A:282:ASP:N	1:A:282:ASP:OD1	2.52	0.42
1:B:14:HIS:CE1	1:B:33:VAL:HG11	2.54	0.42
1:B:470:GLN:OE1	1:B:471:CYS:N	2.51	0.42
1:A:125:THR:O	1:A:153:SER:HB3	2.19	0.42
1:C:282:ASP:OD1	1:C:282:ASP:N	2.53	0.42
1:A:472:ASP:HB3	1:A:474:SER:N	2.31	0.42
1:B:385:LYS:HB3	1:B:386:PRO:CD	2.39	0.42
1:B:221:ARG:HD2	3:B:708:HOH:O	2.19	0.42
1:B:386:PRO:O	1:B:387:ASN:HB3	2.19	0.42
1:B:11:LEU:HD22	1:B:445:LEU:HG	2.02	0.42
1:B:111:ASN:HA	1:B:261:LYS:HD3	2.02	0.42
1:C:454:ARG:HG3	1:C:486:HIS:CG	2.55	0.42
1:B:38:LEU:HD12	1:B:427:VAL:HG22	2.01	0.41
1:B:144:PHE:HB2	1:B:147:LEU:HB2	2.02	0.41
1:B:477:GLU:HG3	1:B:481:ASN:ND2	2.35	0.41
1:A:13:HIS:HB2	1:A:318:MET:SD	2.61	0.41
1:C:125:THR:O	1:C:125:THR:CG2	2.66	0.41
1:C:125:THR:O	1:C:153:SER:HB3	2.20	0.41
1:C:170:ASP:OD1	3:C:710:HOH:O	2.22	0.41
1:C:385:LYS:O	1:C:385:LYS:HE3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:ASN:ND2	1:A:96:PHE:CE1	2.89	0.41
1:A:472:ASP:O	1:A:476:ILE:HG12	2.20	0.41
1:C:385:LYS:CD	3:C:703:HOH:O	2.68	0.40
1:B:34:THR:OG1	1:B:316:THR:O	2.39	0.40
1:C:378:LYS:O	1:C:382:LEU:HD23	2.22	0.40
1:B:380:ASN:CG	3:B:715:HOH:O	2.64	0.40

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	A	482/503 (96%)	463 (96%)	17 (4%)	2 (0%)	30	38
1	B	482/503 (96%)	454 (94%)	25 (5%)	3 (1%)	21	27
1	C	480/503 (95%)	448 (93%)	23 (5%)	9 (2%)	6	5
All	All	1444/1509 (96%)	1365 (94%)	65 (4%)	14 (1%)	12	15

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	472	ASP
1	B	386	PRO
1	C	6	ASN
1	C	338	GLU
1	C	339	ASN
1	C	386	PRO
1	C	387	ASN
1	B	387	ASN
1	C	138	ALA
1	C	125	THR

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Mol	Chain	Res	Type
1	B	6	ASN
1	C	137	ARG
1	C	383	ILE

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	425/437 (97%)	420 (99%)	5 (1%)	63	79
1	B	425/437 (97%)	409 (96%)	16 (4%)	29	44
1	C	424/437 (97%)	410 (97%)	14 (3%)	33	50
All	All	1274/1311 (97%)	1239 (97%)	35 (3%)	39	58

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	151	THR
1	A	259	SER
1	A	260	GLN
1	A	492	GLU
1	A	495	ASN
1	B	17	SER
1	B	90	VAL
1	B	147	LEU
1	B	160	LEU
1	B	163	LEU
1	B	204	GLN
1	B	239	LEU
1	B	261	LYS
1	B	333	ILE
1	B	337	ILE
1	B	342	GLN
1	B	365	LEU
1	B	366	LYS
1	B	404	ILE

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Mol	Chain	Res	Type
1	B	444	LYS
1	B	470	GLN
1	C	9	ILE
1	C	14	HIS
1	C	30	ILE
1	C	137	ARG
1	C	239	LEU
1	C	321	ILE
1	C	336	PHE
1	C	337	ILE
1	C	345	ILE
1	C	357	GLU
1	C	383	ILE
1	C	384	GLU
1	C	477	GLU
1	C	499	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	124	ASN
1	A	204	GLN
1	A	339	ASN
1	A	353	HIS
1	A	495	ASN
1	A	499	GLN
1	B	14	HIS
1	B	42	GLN
1	B	58	GLN
1	B	64	ASN
1	B	339	ASN
1	B	342	GLN
1	B	380	ASN
1	B	387	ASN
1	B	443	ASN
1	C	14	HIS
1	C	74	HIS
1	C	124	ASN
1	C	139	ASN
1	C	339	ASN
1	C	380	ASN

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Mol	Chain	Res	Type
1	C	432	GLN
1	C	481	ASN
1	C	495	ASN
1	C	499	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](#)

6 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	0.52	0	17,19,21	0.51	0
2	NAG	C	601	1	14,14,15	0.32	0	17,19,21	0.59	0
2	NAG	C	602	1	14,14,15	0.69	0	17,19,21	0.47	0
2	NAG	A	602	1	14,14,15	0.47	0	17,19,21	0.45	0
2	NAG	B	601	1	14,14,15	0.25	0	17,19,21	0.44	0
2	NAG	B	602	1	14,14,15	0.52	0	17,19,21	0.53	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	-	2/6/23/26	0/1/1/1
2	NAG	C	601	1	-	2/6/23/26	0/1/1/1
2	NAG	C	602	1	-	0/6/23/26	0/1/1/1
2	NAG	A	602	1	-	2/6/23/26	0/1/1/1
2	NAG	B	601	1	-	0/6/23/26	0/1/1/1
2	NAG	B	602	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	NAG	O5-C5-C6-O6
2	A	602	NAG	O5-C5-C6-O6
2	C	601	NAG	C8-C7-N2-C2
2	C	601	NAG	O7-C7-N2-C2
2	A	601	NAG	C4-C5-C6-O6
2	A	602	NAG	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	601	NAG	1	0
2	C	602	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

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, 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	486/503 (96%)	-0.01	15 (3%) 51 53	15, 30, 64, 105	0
1	B	486/503 (96%)	0.29	33 (6%) 23 25	16, 32, 78, 146	0
1	C	484/503 (96%)	0.54	69 (14%) 6 7	19, 37, 97, 147	0
All	All	1456/1509 (96%)	0.27	117 (8%) 18 20	15, 34, 83, 147	0

All (117) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	336	PHE	8.7
1	B	387	ASN	7.8
1	B	322	PRO	7.0
1	C	337	ILE	6.0
1	B	383	ILE	5.9
1	C	383	ILE	5.8
1	C	335	GLY	5.4
1	C	386	PRO	5.4
1	B	323	GLU	5.3
1	B	389	LYS	5.3
1	B	333	ILE	5.0
1	B	475	CYS	5.0
1	C	356	ALA	4.9
1	B	386	PRO	4.7
1	C	352	ARG	4.6
1	B	471	CYS	4.4
1	C	464	CYS	4.4
1	C	385	LYS	4.3
1	C	351	PHE	4.2
1	B	390	TYR	4.2
1	C	5	GLY	4.0
1	C	323	GLU	3.9
1	C	9	ILE	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	5	GLY	3.6
1	A	499	GLN	3.6
1	C	465	PHE	3.6
1	C	366	LYS	3.5
1	C	345	ILE	3.5
1	C	488	ILE	3.5
1	C	138	ALA	3.4
1	A	64	ASN	3.4
1	B	362	ALA	3.3
1	B	382	LEU	3.3
1	B	358	GLY	3.3
1	B	477	GLU	3.3
1	C	137	ARG	3.2
1	C	474	SER	3.2
1	B	359	THR	3.2
1	B	498	PHE	3.2
1	C	472	ASP	3.1
1	C	321	ILE	3.1
1	C	471	CYS	3.1
1	C	350	GLY	3.1
1	C	8	VAL	3.1
1	B	474	SER	3.1
1	C	11	LEU	3.1
1	C	478	SER	3.0
1	C	353	HIS	3.0
1	C	468	PHE	3.0
1	C	387	ASN	3.0
1	B	384	GLU	3.0
1	C	499	GLN	2.9
1	C	14	HIS	2.9
1	B	470	GLN	2.9
1	C	479	ILE	2.9
1	B	385	LYS	2.8
1	C	342	GLN	2.8
1	B	365	LEU	2.8
1	C	486	HIS	2.8
1	C	384	GLU	2.8
1	C	447	GLU	2.7
1	C	354	GLN	2.7
1	C	482	GLY	2.7
1	C	469	HIS	2.7
1	C	498	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	359	THR	2.7
1	C	362	ALA	2.7
1	C	489	TYR	2.7
1	C	358	GLY	2.7
1	B	388	GLU	2.6
1	B	337	ILE	2.6
1	A	471	CYS	2.6
1	C	494	ILE	2.6
1	A	6	ASN	2.6
1	B	5	GLY	2.6
1	A	356	ALA	2.5
1	B	356	ALA	2.5
1	B	321	ILE	2.5
1	C	339	ASN	2.5
1	C	483	THR	2.5
1	A	498	PHE	2.5
1	B	338	GLU	2.5
1	B	488	ILE	2.5
1	C	470	GLN	2.4
1	A	322	PRO	2.4
1	C	7	PRO	2.4
1	C	338	GLU	2.4
1	A	207	GLN	2.4
1	A	472	ASP	2.3
1	B	380	ASN	2.3
1	C	347	GLY	2.3
1	C	10	CYS	2.3
1	C	473	ASN	2.3
1	B	499	GLN	2.3
1	C	6	ASN	2.3
1	A	14	HIS	2.3
1	B	339	ASN	2.3
1	C	139	ASN	2.3
1	C	360	GLY	2.2
1	C	453	LEU	2.2
1	C	388	GLU	2.2
1	C	348	TRP	2.2
1	C	155	GLY	2.2
1	B	457	ALA	2.2
1	A	386	PRO	2.1
1	A	383	ILE	2.1
1	A	385	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	365	LEU	2.1
1	C	341	TRP	2.1
1	C	477	GLU	2.1
1	C	28	ASP	2.1
1	C	467	ILE	2.1
1	A	495	ASN	2.0
1	C	475	CYS	2.0
1	B	18	ASN	2.0
1	C	382	LEU	2.0
1	C	322	PRO	2.0

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($\text{\AA}^2$)	Q<0.9
2	NAG	C	602	14/15	0.79	0.15	42,48,57,58	0
2	NAG	B	601	14/15	0.80	0.14	68,73,77,80	0
2	NAG	A	601	14/15	0.84	0.13	39,43,51,53	0
2	NAG	A	602	14/15	0.84	0.12	31,37,40,45	0
2	NAG	B	602	14/15	0.88	0.14	38,45,54,54	0
2	NAG	C	601	14/15	0.90	0.10	34,40,46,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.