



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:31 PM UTC

PDB ID : 5G47 / pdb_00005g47
Title : Structure of Gc glycoprotein from severe fever with thrombocytopenia syndrome virus in the trimeric postfusion conformation
Authors : Halldorsson, S.; Behrens, A.J.; Harlos, K.; Huiskonen, J.T.; Elliott, R.M.; Crispin, M.; Brennan, B.; Bowden, T.A.
Deposited on : 2016-05-05
Resolution : 2.45 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

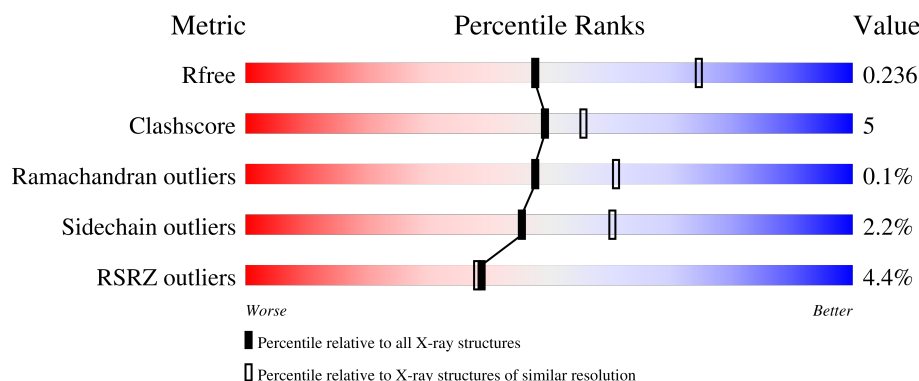
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	445	4% 81% 15% .
1	B	445	4% 77% 19% .
1	C	445	5% 85% 10% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9726 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 SFTSV GC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	428	Total	C	N	O	S	0	0	0
			3194	1982	562	616	34			
1	B	428	Total	C	N	O	S	0	0	0
			3194	1982	562	616	34			
1	C	428	Total	C	N	O	S	0	0	0
			3194	1982	562	616	34			

There are 33 discrepancies between the modelled and reference sequences:

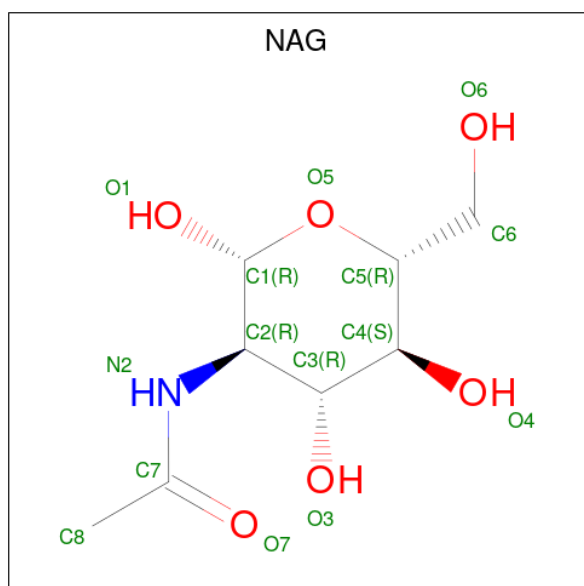
Chain	Residue	Modelled	Actual	Comment	Reference
A	561	THR	-	expression tag	UNP R4V2Q5
A	562	GLY	-	expression tag	UNP R4V2Q5
A	997	GLY	-	expression tag	UNP R4V2Q5
A	998	THR	-	expression tag	UNP R4V2Q5
A	999	LYS	-	expression tag	UNP R4V2Q5
A	1000	HIS	-	expression tag	UNP R4V2Q5
A	1001	HIS	-	expression tag	UNP R4V2Q5
A	1002	HIS	-	expression tag	UNP R4V2Q5
A	1003	HIS	-	expression tag	UNP R4V2Q5
A	1004	HIS	-	expression tag	UNP R4V2Q5
A	1005	HIS	-	expression tag	UNP R4V2Q5
B	561	THR	-	expression tag	UNP R4V2Q5
B	562	GLY	-	expression tag	UNP R4V2Q5
B	997	GLY	-	expression tag	UNP R4V2Q5
B	998	THR	-	expression tag	UNP R4V2Q5
B	999	LYS	-	expression tag	UNP R4V2Q5
B	1000	HIS	-	expression tag	UNP R4V2Q5
B	1001	HIS	-	expression tag	UNP R4V2Q5
B	1002	HIS	-	expression tag	UNP R4V2Q5
B	1003	HIS	-	expression tag	UNP R4V2Q5
B	1004	HIS	-	expression tag	UNP R4V2Q5
B	1005	HIS	-	expression tag	UNP R4V2Q5
C	561	THR	-	expression tag	UNP R4V2Q5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	562	GLY	-	expression tag	UNP R4V2Q5
C	997	GLY	-	expression tag	UNP R4V2Q5
C	998	THR	-	expression tag	UNP R4V2Q5
C	999	LYS	-	expression tag	UNP R4V2Q5
C	1000	HIS	-	expression tag	UNP R4V2Q5
C	1001	HIS	-	expression tag	UNP R4V2Q5
C	1002	HIS	-	expression tag	UNP R4V2Q5
C	1003	HIS	-	expression tag	UNP R4V2Q5
C	1004	HIS	-	expression tag	UNP R4V2Q5
C	1005	HIS	-	expression tag	UNP R4V2Q5

- 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	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		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Cl 1	0	0

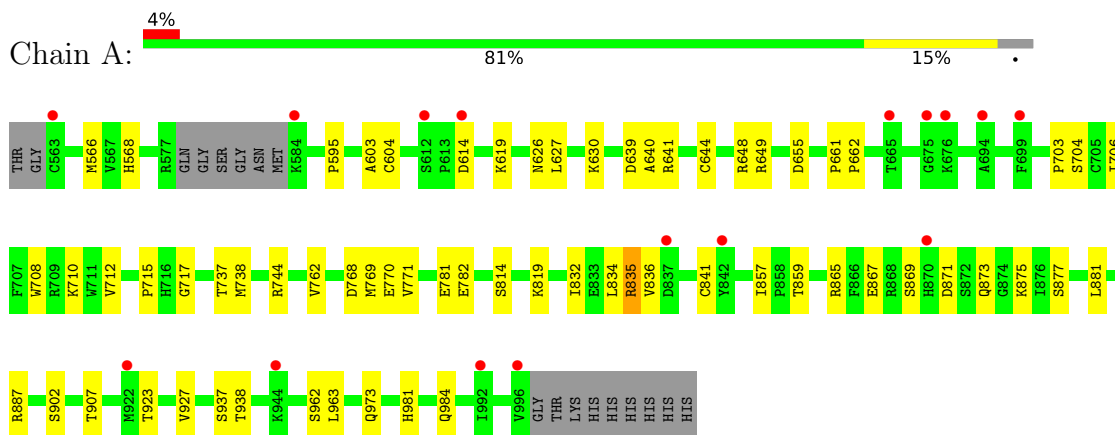
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	42	Total 42	O 42	0	0
4	B	32	Total 32	O 32	0	0
4	C	27	Total 27	O 27	0	0

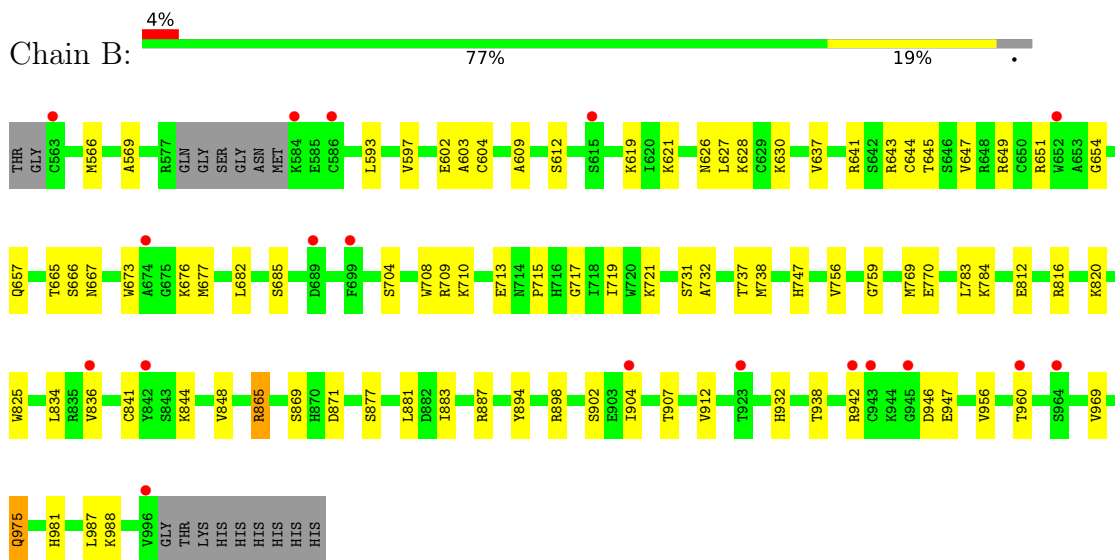
3 Residue-property plots [i](#)

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.

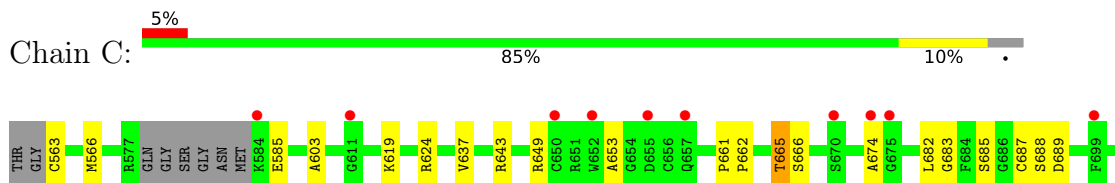
• Molecule 1: SFTSV GC

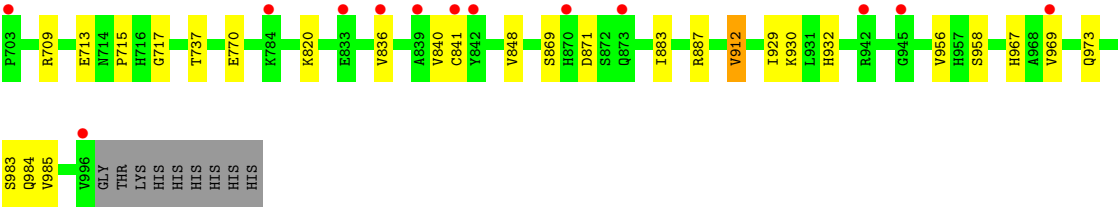


• Molecule 1: SFTSV GC



• Molecule 1: SFTSV GC





4 Data and refinement statistics

Property	Value	Source
Space group	I 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	147.36Å 152.27Å 160.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	108.51 – 2.45 108.51 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.6 (108.51-2.45) 99.6 (108.51-2.45)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.45Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.194 , 0.232 0.199 , 0.236	Depositor DCC
R_{free} test set	3102 reflections (4.68%)	wwPDB-VP
Wilson B-factor (Å ²)	60.6	Xtriage
Anisotropy	0.361	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 49.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.013 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9726	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% 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, CL

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	0.10	0/3255	0.32	0/4401
1	B	0.10	0/3255	0.30	0/4401
1	C	0.11	0/3255	0.33	0/4401
All	All	0.10	0/9765	0.32	0/13203

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

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	3194	0	3123	36	0
1	B	3194	0	3123	47	0
1	C	3194	0	3123	23	0
2	A	14	0	13	0	0
2	B	14	0	13	0	0
2	C	14	0	13	0	0
3	B	1	0	0	0	0
4	A	42	0	0	0	0
4	B	32	0	0	0	0
4	C	27	0	0	2	0
All	All	9726	0	9408	102	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 (102) 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:A:738:MET:HE2	1:A:744:ARG:HD3	1.68	0.74
1:B:619:LYS:HB2	1:B:737:THR:HB	1.68	0.74
1:B:942:ARG:HB2	1:B:975:GLN:HG3	1.73	0.71
1:C:869:SER:OG	1:C:871:ASP:OD1	2.10	0.69
1:B:628:LYS:NZ	1:B:871:ASP:OD1	2.27	0.68
1:B:628:LYS:NZ	1:B:877:SER:OG	2.25	0.66
1:B:738:MET:HE1	1:B:756:VAL:HB	1.78	0.66
1:C:624:ARG:NH2	4:C:2008:HOH:O	2.29	0.64
1:A:769:MET:HE2	1:A:771:VAL:HB	1.80	0.64
1:A:835:ARG:NH2	1:C:687:CYS:O	2.30	0.64
1:A:973:GLN:OE1	1:A:984:GLN:NE2	2.33	0.62
1:A:770:GLU:OE2	1:A:887:ARG:NH2	2.33	0.61
1:C:563:CYS:N	4:C:2001:HOH:O	2.33	0.61
1:A:781:GLU:OE2	1:A:875:LYS:NZ	2.31	0.60
1:A:619:LYS:HB2	1:A:737:THR:HB	1.82	0.60
1:A:865:ARG:NH1	1:A:867:GLU:OE1	2.32	0.60
1:B:609:ALA:HB3	1:B:612:SER:HB3	1.86	0.57
1:C:619:LYS:HB2	1:C:737:THR:HB	1.85	0.56
1:B:641:ARG:NH2	1:B:713:GLU:OE2	2.37	0.56
1:B:987:LEU:O	1:B:988:LYS:NZ	2.37	0.56
1:B:667:ASN:OD1	1:B:820:LYS:NZ	2.38	0.56
1:B:869:SER:OG	1:B:871:ASP:OD1	2.24	0.56
1:B:942:ARG:HG2	1:B:947:GLU:HG3	1.88	0.55
1:A:869:SER:OG	1:A:871:ASP:OD1	2.23	0.54
1:B:654:GLY:O	1:B:657:GLN:NE2	2.40	0.54
1:C:643:ARG:NH1	1:C:713:GLU:OE2	2.41	0.53
1:B:770:GLU:OE2	1:B:887:ARG:NH2	2.42	0.53
1:B:732:ALA:O	1:B:747:HIS:ND1	2.24	0.53
1:B:717:GLY:O	1:B:816:ARG:NH1	2.42	0.52
1:B:645:THR:HG21	1:B:677:MET:HE3	1.90	0.52
1:A:626:ASN:HB3	1:A:877:SER:HB2	1.91	0.52
1:A:927:VAL:HG21	1:A:963:LEU:HD12	1.91	0.52
1:C:683:GLY:O	1:C:820:LYS:NZ	2.30	0.50
1:A:604:CYS:HB2	1:A:902:SER:HB3	1.94	0.49
1:A:881:LEU:HD13	1:A:962:SER:HB3	1.94	0.49
1:B:627:LEU:HD11	1:B:769:MET:SD	2.53	0.49
1:B:626:ASN:O	1:B:731:SER:N	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:644:CYS:HB3	1:B:710:LYS:HA	1.95	0.48
1:B:602:GLU:HG3	1:B:621:LYS:HB2	1.95	0.48
1:A:644:CYS:HB3	1:A:710:LYS:HA	1.96	0.48
1:A:708:TRP:CD1	1:A:834:LEU:HD11	2.49	0.47
1:C:682:LEU:HD12	1:C:820:LYS:HD3	1.95	0.47
1:B:932:HIS:HB2	1:B:956:VAL:HG22	1.96	0.47
1:C:930:LYS:HA	1:C:958:SER:HA	1.97	0.47
1:A:566:MET:HB2	1:A:603:ALA:HB2	1.95	0.47
1:A:627:LEU:HD11	1:A:769:MET:CE	2.45	0.47
1:B:715:PRO:C	1:B:717:GLY:H	2.21	0.47
1:A:710:LYS:HD3	1:A:832:ILE:HG22	1.97	0.47
1:A:627:LEU:HD11	1:A:769:MET:HE1	1.96	0.47
1:B:676:LYS:HD2	1:B:676:LYS:HA	1.66	0.46
1:A:703:PRO:HG2	1:B:836:VAL:HG21	1.97	0.46
1:A:873:GLN:HE22	1:B:946:ASP:HA	1.81	0.46
1:B:569:ALA:HB3	1:B:593:LEU:HB3	1.97	0.46
1:C:649:ARG:HH21	1:C:653:ALA:HB1	1.80	0.46
1:A:715:PRO:C	1:A:717:GLY:H	2.24	0.46
1:B:865:ARG:HB2	1:B:881:LEU:HD21	1.98	0.46
1:B:566:MET:HB2	1:B:603:ALA:HB2	1.98	0.45
1:B:665:THR:OG1	1:B:666:SER:N	2.47	0.45
1:C:637:VAL:HG11	1:C:848:VAL:HG21	1.98	0.45
1:B:883:ILE:HD11	1:B:938:THR:HG23	1.99	0.45
1:C:566:MET:HB2	1:C:603:ALA:HB2	1.99	0.45
1:A:639:ASP:CG	1:A:641:ARG:HH12	2.26	0.44
1:A:648:ARG:HB2	1:A:706:ILE:HD13	1.99	0.44
1:A:640:ALA:HB1	1:A:712:VAL:CG2	2.47	0.44
1:A:907:THR:HG22	1:A:981:HIS:CG	2.52	0.44
1:B:566:MET:HE2	1:B:597:VAL:HG13	1.98	0.44
1:C:770:GLU:OE2	1:C:887:ARG:NH2	2.51	0.44
1:A:814:SER:HA	1:A:819:LYS:HE2	2.00	0.44
1:B:759:GLY:O	1:B:894:TYR:HA	2.17	0.44
1:C:715:PRO:C	1:C:717:GLY:H	2.26	0.44
1:A:649:ARG:O	1:A:704:SER:HB2	2.18	0.43
1:C:683:GLY:H	1:C:820:LYS:HE2	1.83	0.43
1:C:912:VAL:HG12	1:C:985:VAL:HB	2.00	0.43
1:B:685:SER:OG	1:B:709:ARG:NH1	2.44	0.43
1:C:983:SER:OG	1:C:984:GLN:N	2.52	0.43
1:B:630:LYS:HA	1:B:630:LYS:HD2	1.80	0.43
1:C:665:THR:OG1	1:C:666:SER:N	2.46	0.43
1:A:568:HIS:HD2	1:A:595:PRO:HD2	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:604:CYS:HB2	1:B:902:SER:HB3	1.99	0.43
1:B:721:LYS:HB2	1:B:783:LEU:HD21	2.00	0.43
1:B:651:ARG:HA	1:B:651:ARG:HD3	1.91	0.42
1:B:708:TRP:CD1	1:B:834:LEU:HD11	2.54	0.42
1:C:685:SER:HA	1:C:709:ARG:HA	2.00	0.42
1:C:967:HIS:HD2	1:C:969:VAL:O	2.02	0.42
1:B:641:ARG:HH12	1:B:643:ARG:HD3	1.85	0.42
1:A:655:ASP:OD1	1:A:661:PRO:HG2	2.20	0.41
1:C:661:PRO:HA	1:C:662:PRO:HD3	1.94	0.41
1:B:637:VAL:HG11	1:B:848:VAL:HG21	2.01	0.41
1:B:647:VAL:HG13	1:B:673:TRP:HD1	1.85	0.41
1:B:784:LYS:NZ	1:B:812:GLU:OE2	2.51	0.41
1:C:932:HIS:HB2	1:C:956:VAL:HG22	2.03	0.41
1:A:630:LYS:HD2	1:A:630:LYS:HA	1.84	0.41
1:A:836:VAL:N	1:C:689:ASP:OD2	2.45	0.41
1:A:857:ILE:HD13	1:A:859:THR:HG22	2.03	0.41
1:A:937:SER:OG	1:A:938:THR:N	2.54	0.41
1:B:649:ARG:O	1:B:704:SER:HB2	2.21	0.41
1:B:682:LEU:HD21	1:B:825:TRP:CG	2.56	0.41
1:A:768:ASP:HB2	1:A:887:ARG:HG2	2.03	0.40
1:A:661:PRO:HA	1:A:662:PRO:HD3	1.95	0.40
1:B:719:ILE:HD11	1:B:816:ARG:HA	2.04	0.40
1:B:907:THR:HG22	1:B:981:HIS:CG	2.56	0.40
1:B:912:VAL:HG11	1:B:987:LEU:HD23	2.02	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	424/445 (95%)	408 (96%)	16 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	424/445 (95%)	411 (97%)	13 (3%)	0	100	100
1	C	424/445 (95%)	413 (97%)	10 (2%)	1 (0%)	43	54
All	All	1272/1335 (95%)	1232 (97%)	39 (3%)	1 (0%)	48	61

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	674	ALA

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	358/371 (96%)	352 (98%)	6 (2%)	53	67
1	B	358/371 (96%)	350 (98%)	8 (2%)	45	60
1	C	358/371 (96%)	348 (97%)	10 (3%)	38	54
All	All	1074/1113 (96%)	1050 (98%)	24 (2%)	45	60

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	614	ASP
1	A	762	VAL
1	A	782	GLU
1	A	835	ARG
1	A	841	CYS
1	A	923	THR
1	B	841	CYS
1	B	844	LYS
1	B	865	ARG
1	B	898	ARG
1	B	904	ILE
1	B	960	THR
1	B	969	VAL

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Mol	Chain	Res	Type
1	B	975	GLN
1	C	585	GLU
1	C	665	THR
1	C	688	SER
1	C	836	VAL
1	C	840	VAL
1	C	841	CYS
1	C	883	ILE
1	C	912	VAL
1	C	929	ILE
1	C	973	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	598	ASN
1	A	973	GLN
1	A	975	GLN
1	A	984	GLN
1	A	990	ASN
1	B	716	HIS
1	C	606	HIS
1	C	755	GLN
1	C	932	HIS
1	C	973	GLN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 for Mogul analysis.

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	C	1997	1	14,14,15	0.17	0	17,19,21	0.52	0
2	NAG	B	1997	1	14,14,15	0.20	0	17,19,21	0.57	0
2	NAG	A	1997	1	14,14,15	0.21	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
2	NAG	C	1997	1	-	4/6/23/26	0/1/1/1
2	NAG	B	1997	1	-	4/6/23/26	0/1/1/1
2	NAG	A	1997	1	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1997	NAG	O5-C5-C6-O6
2	A	1997	NAG	C4-C5-C6-O6
2	A	1997	NAG	C8-C7-N2-C2
2	A	1997	NAG	O7-C7-N2-C2
2	B	1997	NAG	C8-C7-N2-C2
2	B	1997	NAG	O7-C7-N2-C2
2	C	1997	NAG	C8-C7-N2-C2
2	C	1997	NAG	O7-C7-N2-C2
2	B	1997	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
2	B	1997	NAG	C4-C5-C6-O6
2	C	1997	NAG	C4-C5-C6-O6
2	C	1997	NAG	O5-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, 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	428/445 (96%)	0.32	16 (3%) 45 45	43, 72, 132, 164	0
1	B	428/445 (96%)	0.37	18 (4%) 40 40	41, 81, 138, 181	0
1	C	428/445 (96%)	0.35	23 (5%) 31 29	42, 74, 124, 153	0
All	All	1284/1335 (96%)	0.35	57 (4%) 39 38	41, 75, 132, 181	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	996	VAL	5.2
1	B	945	GLY	5.2
1	C	674	ALA	5.1
1	C	842	TYR	4.4
1	C	996	VAL	4.2
1	A	842	TYR	4.1
1	B	996	VAL	3.3
1	A	563	CYS	3.3
1	B	584	LYS	3.3
1	B	904	ILE	3.3
1	C	699	PHE	3.1
1	B	842	TYR	3.0
1	A	675	GLY	2.9
1	C	945	GLY	2.9
1	C	839	ALA	2.9
1	A	699	PHE	2.9
1	C	836	VAL	2.8
1	C	833	GLU	2.8
1	C	584	LYS	2.8
1	C	873	GLN	2.7
1	B	652	TRP	2.6
1	A	584	LYS	2.6
1	B	923	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	964	SER	2.5
1	A	665	THR	2.5
1	C	650	CYS	2.5
1	C	942	ARG	2.5
1	B	586	CYS	2.4
1	A	676	LYS	2.4
1	B	943	CYS	2.4
1	C	784	LYS	2.4
1	A	612	SER	2.4
1	B	563	CYS	2.4
1	C	652	TRP	2.3
1	A	870	HIS	2.3
1	A	922	MET	2.3
1	A	837	ASP	2.2
1	C	657	GLN	2.2
1	C	670	SER	2.2
1	A	614	ASP	2.2
1	C	969	VAL	2.1
1	C	675	GLY	2.1
1	C	870	HIS	2.1
1	C	611	GLY	2.1
1	C	655	ASP	2.1
1	B	615	SER	2.1
1	C	841	CYS	2.1
1	B	836	VAL	2.1
1	A	992	ILE	2.1
1	B	689	ASP	2.1
1	B	699	PHE	2.0
1	A	694	ALA	2.0
1	B	942	ARG	2.0
1	B	960	THR	2.0
1	C	703	PRO	2.0
1	B	674	ALA	2.0
1	A	944	LYS	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 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	1997	14/15	0.76	0.14	89,112,136,137	0
2	NAG	B	1997	14/15	0.83	0.13	79,94,105,108	0
2	NAG	A	1997	14/15	0.84	0.13	79,100,109,122	0
3	CL	B	1998	1/1	0.96	0.21	87,87,87,87	0

6.5 Other polymers [i](#)

There are no such residues in this entry.