



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:49 PM UTC

PDB ID : 6JYO / pdb_00006jyo
Title : GII.13/21 noroviruses recognize glycans with a terminal beta-galactose via an unconventional glycan binding site
Authors : Duan, Z.; Xin, C.
Deposited on : 2019-04-27
Resolution : 1.50 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
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

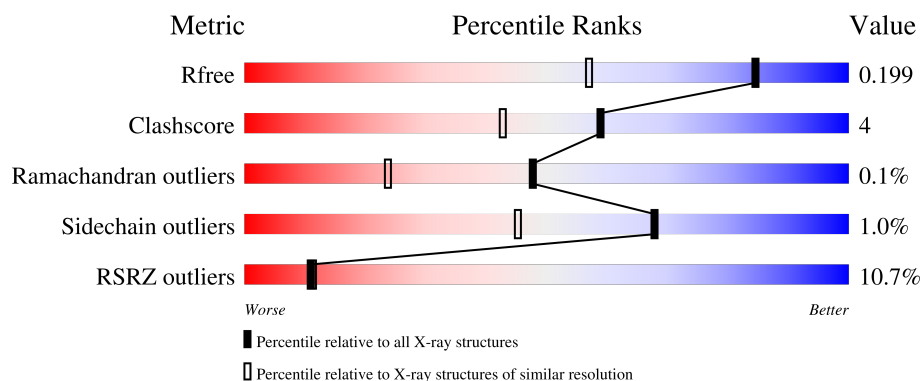
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

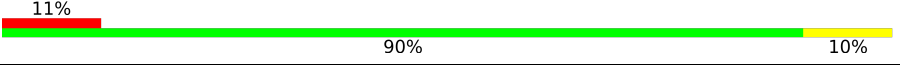
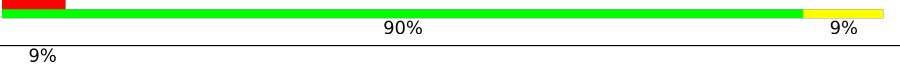
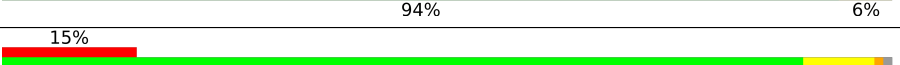
The reported resolution of this entry is 1.50 Å.

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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	309	 11% 90% 10%
1	B	309	 7% 90% 9%
1	C	309	 9% 94% 6%
1	D	309	 15% 90% 8% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11206 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 norovirus P domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	309	Total	C	N	O	S	0	1	0
			2401	1525	410	460	6			
1	B	309	Total	C	N	O	S	0	0	0
			2396	1522	409	459	6			
1	C	309	Total	C	N	O	S	0	0	0
			2396	1522	409	459	6			
1	D	307	Total	C	N	O	S	0	0	0
			2381	1514	407	454	6			

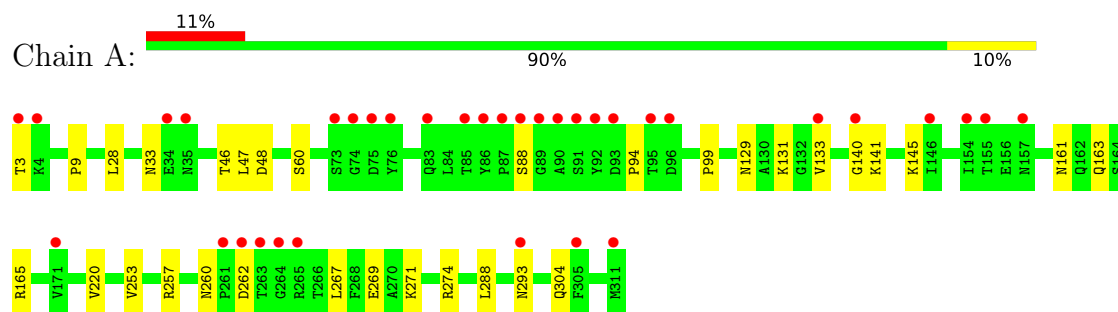
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	364	Total	O	0	0
			364	364		
2	B	449	Total	O	0	0
			449	449		
2	C	416	Total	O	0	0
			416	416		
2	D	403	Total	O	0	0
			403	403		

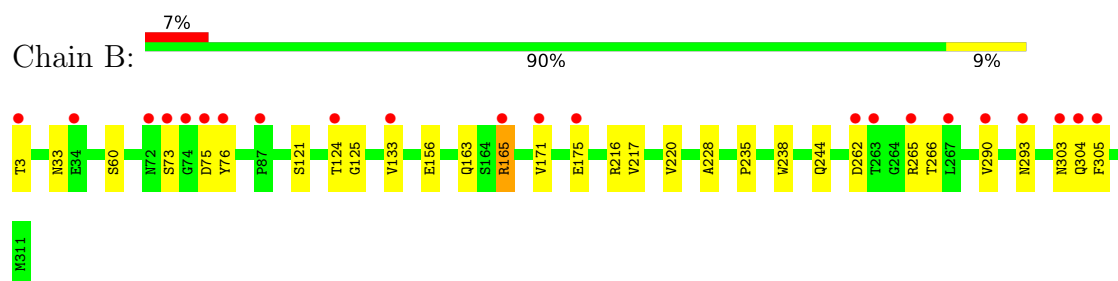
3 Residue-property plots [i](#)

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.

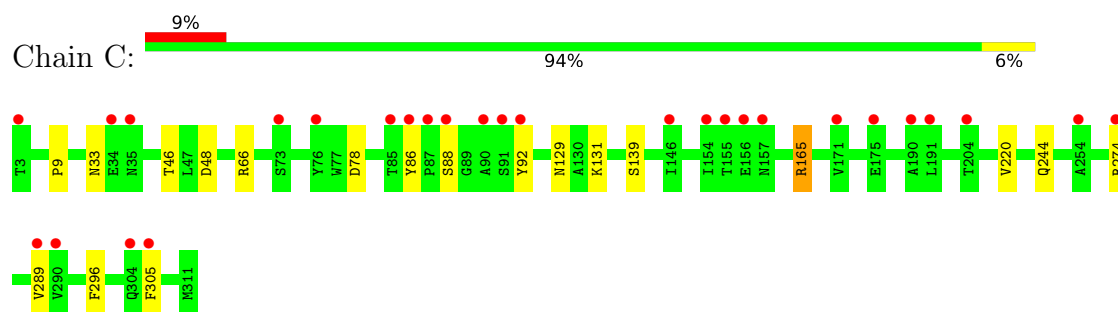
- Molecule 1: norovirus P domain protein



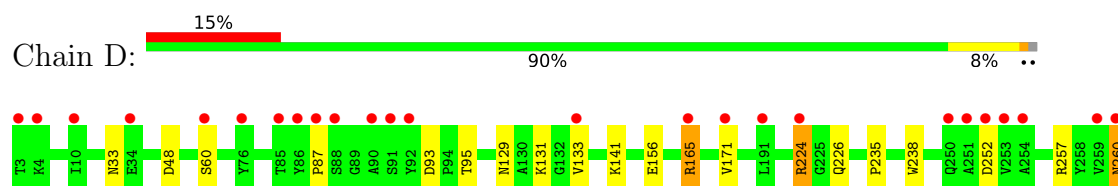
- Molecule 1: norovirus P domain protein

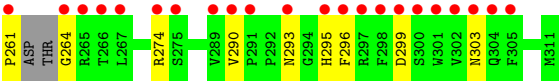


- Molecule 1: norovirus P domain protein



- Molecule 1: norovirus P domain protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	87.64Å 105.67Å 138.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.35 – 1.50 49.35 – 1.50	Depositor EDS
% Data completeness (in resolution range)	98.9 (49.35-1.50) 98.9 (49.35-1.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 1.50Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.188 , 0.199 0.188 , 0.199	Depositor DCC
R_{free} test set	10097 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	14.1	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11206	wwPDB-VP
Average B, all atoms (Å ²)	19.0	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 65.88 % 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 6.6027e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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	0.19	0/2471	0.39	0/3383
1	B	0.19	0/2463	0.45	2/3372 (0.1%)
1	C	0.17	0/2463	0.41	2/3372 (0.1%)
1	D	0.25	1/2447 (0.0%)	0.47	3/3348 (0.1%)
All	All	0.20	1/9844 (0.0%)	0.43	7/13475 (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
1	B	0	1
1	C	0	1
1	D	0	2
All	All	0	5

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	156	GLU	CD-OE2	-5.47	1.15	1.25

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	165	ARG	NE-CZ-NH2	7.47	125.92	119.20
1	B	165	ARG	NE-CZ-NH1	-7.31	114.19	121.50
1	C	165	ARG	NE-CZ-NH1	-6.32	115.18	121.50
1	C	165	ARG	NE-CZ-NH2	6.23	124.80	119.20
1	D	165	ARG	NE-CZ-NH2	6.16	124.74	119.20
1	D	224	ARG	NE-CZ-NH1	-5.22	116.28	121.50
1	D	224	ARG	NE-CZ-NH2	5.03	123.73	119.20

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	33	ASN	Peptide
1	B	33	ASN	Peptide
1	C	33	ASN	Peptide
1	D	260	ASN	Peptide
1	D	33	ASN	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	2401	0	2315	21	1
1	B	2396	0	2309	26	0
1	C	2396	0	2309	14	0
1	D	2381	0	2297	27	1
2	A	364	0	0	10	0
2	B	449	0	0	13	3
2	C	416	0	0	7	2
2	D	403	0	0	13	6
All	All	11206	0	9230	82	9

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 4.

All (82) 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:D:165:ARG:NH1	2:D:401:HOH:O	1.84	1.07
1:D:293:ASN:OD1	2:D:402:HOH:O	1.85	0.93
1:B:165:ARG:NH1	2:B:401:HOH:O	2.00	0.93
1:B:3:THR:N	2:B:402:HOH:O	2.02	0.93
1:B:303:ASN:HA	2:B:585:HOH:O	1.69	0.91
1:D:87:PRO:O	2:D:403:HOH:O	1.90	0.88
1:B:60:SER:HB3	2:B:698:HOH:O	1.79	0.82
1:A:94:PRO:O	1:A:145:LYS:NZ	2.19	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:264:GLY:N	2:D:405:HOH:O	2.22	0.73
1:B:304:GLN:HG2	2:B:648:HOH:O	1.90	0.71
1:D:226:GLN:NE2	2:D:406:HOH:O	2.23	0.71
1:A:274:ARG:NH1	2:A:409:HOH:O	2.23	0.71
1:B:175:GLU:H	1:B:175:GLU:CD	2.00	0.70
1:D:264:GLY:O	2:D:404:HOH:O	2.09	0.69
1:D:165:ARG:HD2	2:D:719:HOH:O	1.92	0.68
1:B:76:TYR:OH	2:B:403:HOH:O	2.07	0.66
1:B:244:GLN:HG2	2:B:685:HOH:O	1.95	0.66
1:D:48:ASP:CG	1:D:274:ARG:HH22	2.05	0.64
1:B:163:GLN:NE2	2:B:406:HOH:O	2.26	0.63
1:C:92:TYR:HA	2:C:406:HOH:O	1.98	0.63
1:D:93:ASP:OD2	1:D:95:THR:OG1	2.11	0.61
1:A:165:ARG:NH2	2:A:407:HOH:O	2.22	0.60
2:B:566:HOH:O	1:D:60:SER:HB3	2.03	0.58
1:D:257:ARG:HH22	1:D:299:ASP:CG	2.11	0.58
1:B:262:ASP:OD2	1:B:293:ASN:ND2	2.36	0.58
1:D:257:ARG:NH2	1:D:299:ASP:OD1	2.35	0.57
1:D:295:HIS:HB2	2:D:410:HOH:O	2.05	0.56
1:A:262:ASP:OD2	1:A:293:ASN:ND2	2.38	0.56
1:B:121:SER:HB3	1:D:224:ARG:CZ	2.37	0.55
1:D:257:ARG:HH22	1:D:299:ASP:HB2	1.70	0.55
1:B:73:SER:HB3	1:C:305:PHE:CZ	2.42	0.54
1:D:303:ASN:ND2	2:D:414:HOH:O	2.40	0.54
1:B:266:THR:OG1	2:B:404:HOH:O	2.10	0.54
1:C:66:ARG:HG2	1:C:165:ARG:HG3	1.91	0.53
1:C:274:ARG:HD2	2:C:723:HOH:O	2.08	0.53
1:A:141:LYS:NZ	2:A:418:HOH:O	2.38	0.53
1:B:303:ASN:OD1	1:B:305:PHE:N	2.39	0.53
1:A:60:SER:O	2:A:404:HOH:O	2.19	0.53
1:D:141:LYS:NZ	2:D:416:HOH:O	2.42	0.52
1:D:260:ASN:O	1:D:261:PRO:O	2.28	0.51
1:D:257:ARG:HH22	1:D:299:ASP:CB	2.24	0.51
1:D:296:PHE:HB2	2:D:447:HOH:O	2.10	0.51
1:B:217:VAL:HG12	1:B:228:ALA:O	2.11	0.51
1:D:299:ASP:HA	2:D:420:HOH:O	2.10	0.51
1:A:304:GLN:HG2	2:A:406:HOH:O	2.11	0.50
1:A:129:ASN:HB3	1:C:220:VAL:HA	1.93	0.50
1:A:140:GLY:HA2	2:A:567:HOH:O	2.13	0.49
1:B:265:ARG:NE	1:B:265:ARG:HA	2.28	0.49
1:B:121:SER:H	1:D:224:ARG:NH2	2.10	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:9:PRO:HB3	2:C:636:HOH:O	2.13	0.48
1:B:220:VAL:HA	1:D:129:ASN:HB3	1.95	0.48
1:A:46:THR:OG1	1:A:48:ASP:OD1	2.29	0.48
1:A:28:LEU:HB2	1:A:288:LEU:HB2	1.96	0.47
1:A:260:ASN:HB2	1:A:267:LEU:HD21	1.96	0.47
1:A:220:VAL:HA	1:C:129:ASN:HB3	1.97	0.47
1:A:88:SER:O	2:A:405:HOH:O	2.20	0.46
1:B:73:SER:C	1:B:75:ASP:H	2.24	0.45
1:B:124:THR:HB	2:B:430:HOH:O	2.17	0.45
1:B:216:ARG:NH1	2:B:423:HOH:O	2.49	0.45
1:D:274:ARG:HH21	1:D:274:ARG:HG2	1.80	0.45
1:A:47:LEU:HG	1:A:274:ARG:HA	1.98	0.45
1:D:235:PRO:HG2	1:D:238:TRP:CD1	2.52	0.45
1:D:257:ARG:NH2	1:D:299:ASP:HB2	2.31	0.45
1:A:9:PRO:HB3	2:A:641:HOH:O	2.17	0.44
1:C:86:TYR:HE2	2:C:406:HOH:O	2.01	0.44
1:D:293:ASN:CG	2:D:402:HOH:O	2.45	0.44
1:A:3:THR:HA	2:A:503:HOH:O	2.17	0.43
1:A:163:GLN:NE2	2:A:423:HOH:O	2.51	0.43
1:A:257:ARG:NH1	1:A:269:GLU:OE1	2.51	0.43
1:B:175:GLU:CD	1:B:175:GLU:N	2.73	0.43
1:C:78:ASP:OD2	1:C:139:SER:OG	2.28	0.43
1:C:88:SER:CB	2:C:428:HOH:O	2.67	0.43
1:C:165:ARG:HH21	1:C:165:ARG:HG2	1.83	0.43
1:B:235:PRO:HG2	1:B:238:TRP:CD1	2.54	0.42
1:C:46:THR:OG1	1:C:48:ASP:OD1	2.36	0.42
1:A:99:PRO:O	1:A:145:LYS:HD3	2.20	0.42
1:B:304:GLN:HG3	1:B:305:PHE:CD2	2.55	0.42
1:C:296:PHE:HB2	2:C:558:HOH:O	2.20	0.41
1:A:253:VAL:HG11	1:A:271:LYS:HG2	2.03	0.41
1:B:125:GLY:N	2:B:430:HOH:O	2.53	0.41
1:C:244:GLN:HG2	2:C:628:HOH:O	2.21	0.40
1:B:303:ASN:OD1	1:B:305:PHE:HB2	2.21	0.40

All (9) 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)
2:D:514:HOH:O	2:D:514:HOH:O[2_665]	1.35	0.85
2:B:834:HOH:O	2:B:841:HOH:O[4_455]	2.06	0.14

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:740:HOH:O	2:D:677:HOH:O[2_655]	2.12	0.08
2:B:809:HOH:O	2:B:810:HOH:O[4_455]	2.15	0.05
1:A:161[B]:ASN:ND2	1:D:252:ASP:OD2[1_455]	2.16	0.04
2:C:799:HOH:O	2:D:791:HOH:O[1_455]	2.17	0.03
2:D:535:HOH:O	2:D:725:HOH:O[2_655]	2.18	0.02
2:C:495:HOH:O	2:D:708:HOH:O[2_665]	2.19	0.01
2:D:404:HOH:O	2:D:514:HOH:O[2_665]	2.19	0.01

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	308/309 (100%)	298 (97%)	10 (3%)	0	100	100
1	B	307/309 (99%)	294 (96%)	12 (4%)	1 (0%)	36	17
1	C	307/309 (99%)	299 (97%)	8 (3%)	0	100	100
1	D	303/309 (98%)	288 (95%)	15 (5%)	0	100	100
All	All	1225/1236 (99%)	1179 (96%)	45 (4%)	1 (0%)	48	24

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	156	GLU

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	264/263 (100%)	262 (99%)	2 (1%)	73	54
1	B	263/263 (100%)	260 (99%)	3 (1%)	65	41
1	C	263/263 (100%)	261 (99%)	2 (1%)	73	54
1	D	261/263 (99%)	257 (98%)	4 (2%)	57	31
All	All	1051/1052 (100%)	1040 (99%)	11 (1%)	68	45

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	131	LYS
1	A	133	VAL
1	B	133	VAL
1	B	171	VAL
1	B	290	VAL
1	C	131	LYS
1	C	289	VAL
1	D	131	LYS
1	D	133	VAL
1	D	171	VAL
1	D	290	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	GLN
1	A	163	GLN
1	A	236	GLN
1	A	240	ASN
1	A	250	GLN
1	B	105	GLN
1	B	236	GLN
1	B	240	ASN
1	B	304	GLN
1	C	159	HIS
1	C	196	ASN
1	C	236	GLN
1	C	240	ASN
1	D	36	ASN
1	D	105	GLN
1	D	196	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	309/309 (100%)	0.58	35 (11%) 10 10	9, 17, 42, 82	1 (0%)
1	B	309/309 (100%)	0.27	22 (7%) 22 24	7, 14, 30, 63	0
1	C	309/309 (100%)	0.38	28 (9%) 15 15	8, 15, 30, 69	0
1	D	307/309 (99%)	0.68	47 (15%) 5 5	8, 17, 36, 65	0
All	All	1234/1236 (99%)	0.48	132 (10%) 11 11	7, 16, 36, 82	1 (0%)

All (132) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	86	TYR	9.3
1	A	90	ALA	9.0
1	C	76	TYR	8.4
1	C	305	PHE	7.8
1	B	305	PHE	7.3
1	D	264	GLY	7.3
1	C	191	LEU	7.0
1	D	261	PRO	7.0
1	B	74	GLY	6.5
1	D	87	PRO	6.4
1	A	76	TYR	6.3
1	A	89	GLY	6.3
1	A	85	THR	6.1
1	C	190	ALA	6.0
1	A	87	PRO	6.0
1	B	73	SER	5.9
1	A	91	SER	5.7
1	C	87	PRO	5.5
1	B	76	TYR	5.4
1	D	259	VAL	5.4
1	D	260	ASN	5.2

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Mol	Chain	Res	Type	RSRZ
1	A	3	THR	5.1
1	D	3	THR	4.7
1	C	155	THR	4.7
1	D	293	ASN	4.7
1	C	34	GLU	4.7
1	B	124	THR	4.6
1	B	304	GLN	4.6
1	D	250	GLN	4.5
1	A	92	TYR	4.4
1	A	155	THR	4.4
1	D	290	VAL	4.3
1	C	3	THR	4.3
1	A	34	GLU	4.3
1	C	304	GLN	4.2
1	A	305	PHE	4.2
1	D	86	TYR	4.2
1	D	171	VAL	4.2
1	B	34	GLU	4.1
1	D	301	TRP	4.1
1	C	154	ILE	4.1
1	A	88	SER	4.1
1	A	263	THR	4.1
1	D	305	PHE	4.0
1	D	85	THR	4.0
1	D	252	ASP	3.9
1	D	304	GLN	3.9
1	B	303	ASN	3.8
1	D	253	VAL	3.7
1	D	297	ARG	3.7
1	C	88	SER	3.7
1	C	91	SER	3.7
1	D	265	ARG	3.6
1	D	34	GLU	3.6
1	A	73	SER	3.5
1	A	171	VAL	3.5
1	C	85	THR	3.5
1	D	274	ARG	3.5
1	D	267	LEU	3.4
1	C	156	GLU	3.3
1	D	10	ILE	3.3
1	B	171	VAL	3.3
1	C	86	TYR	3.3

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Mol	Chain	Res	Type	RSRZ
1	C	171	VAL	3.3
1	D	302	VAL	3.3
1	B	262	ASP	3.2
1	B	290	VAL	3.2
1	D	224	ARG	3.1
1	C	90	ALA	3.1
1	C	289	VAL	3.1
1	D	88	SER	3.0
1	A	75	ASP	2.9
1	B	75	ASP	2.9
1	B	72	ASN	2.9
1	B	265	ARG	2.9
1	B	3	THR	2.9
1	A	261	PRO	2.9
1	B	133	VAL	2.9
1	B	175	GLU	2.9
1	D	289	VAL	2.8
1	D	254	ALA	2.7
1	A	146	ILE	2.6
1	A	95	THR	2.6
1	B	87	PRO	2.6
1	A	74	GLY	2.6
1	C	204	THR	2.5
1	C	175	GLU	2.5
1	D	298	PHE	2.5
1	D	133	VAL	2.5
1	D	299	ASP	2.5
1	D	191	LEU	2.4
1	D	303	ASN	2.4
1	D	296	PHE	2.4
1	D	90	ALA	2.4
1	D	300	SER	2.4
1	D	291	PRO	2.4
1	C	146	ILE	2.4
1	D	266	THR	2.4
1	A	311	MET	2.4
1	D	275	SER	2.3
1	A	157	ASN	2.3
1	A	293	ASN	2.3
1	A	140	GLY	2.3
1	A	262	ASP	2.3
1	B	165	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	263	THR	2.3
1	C	92	TYR	2.3
1	D	92	TYR	2.3
1	C	35	ASN	2.3
1	A	83	GLN	2.2
1	C	290	VAL	2.2
1	D	4	LYS	2.2
1	B	293	ASN	2.2
1	A	93	ASP	2.2
1	D	251	ALA	2.2
1	D	60	SER	2.2
1	A	133	VAL	2.2
1	D	295	HIS	2.2
1	D	76	TYR	2.1
1	A	265	ARG	2.1
1	C	274	ARG	2.1
1	A	35	ASN	2.1
1	A	264	GLY	2.1
1	B	267	LEU	2.1
1	A	4	LYS	2.0
1	D	165	ARG	2.0
1	A	154	ILE	2.0
1	C	73	SER	2.0
1	C	254	ALA	2.0
1	D	91	SER	2.0
1	A	96	ASP	2.0
1	C	157	ASN	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.