



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:25 PM UTC

PDB ID : 5U15 / pdb_00005u15
Title : Crystal Structure of DH270.UCA3 (unliganded) from the DH270 Broadly Neutralizing N332-glycan Dependent Lineage
Authors : Fera, D.; Harrison, S.C.
Deposited on : 2016-11-27
Resolution : 2.26 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

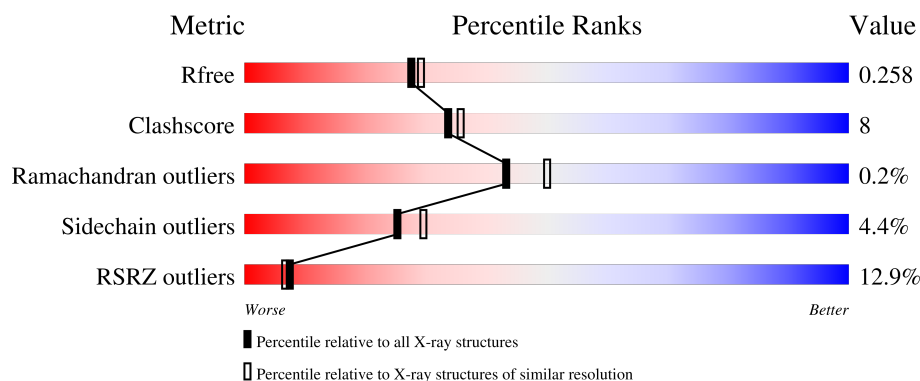
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.26 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	238	
1	H	238	
2	B	216	
2	L	216	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6564 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 DH270.UCA3 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	0
			1579	1000	265	306	8			
1	H	221	Total	C	N	O	S	0	1	0
			1697	1079	285	325	8			

- Molecule 2 is a protein called DH270.UCA3 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	199	Total	C	N	O	S	0	0	0
			1462	915	239	302	6			
2	L	211	Total	C	N	O	S	0	0	0
			1552	968	258	320	6			

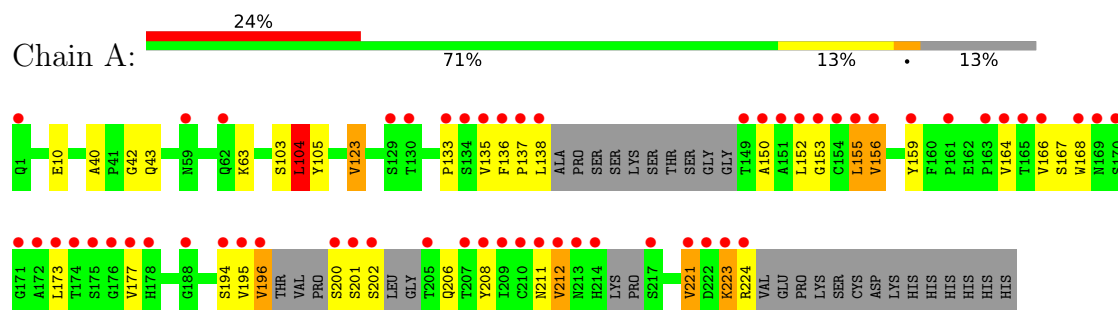
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	70	Total	O	0	0
			70	70		
3	B	42	Total	O	0	0
			42	42		
3	H	89	Total	O	0	0
			89	89		
3	L	73	Total	O	0	0
			73	73		

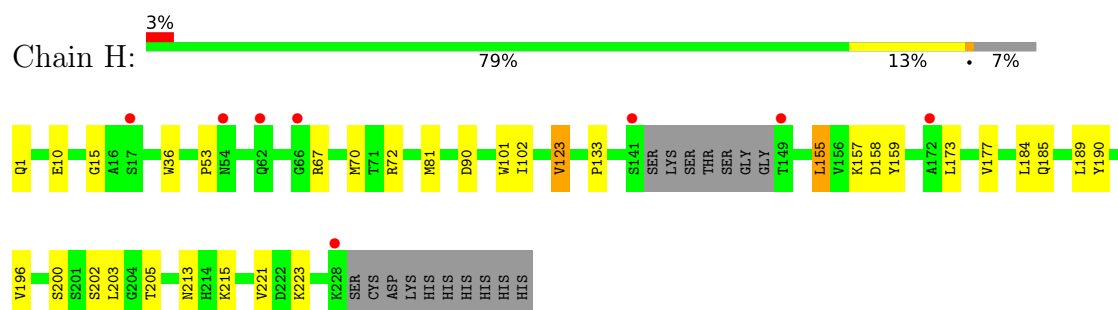
3 Residue-property plots [i](#)

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.

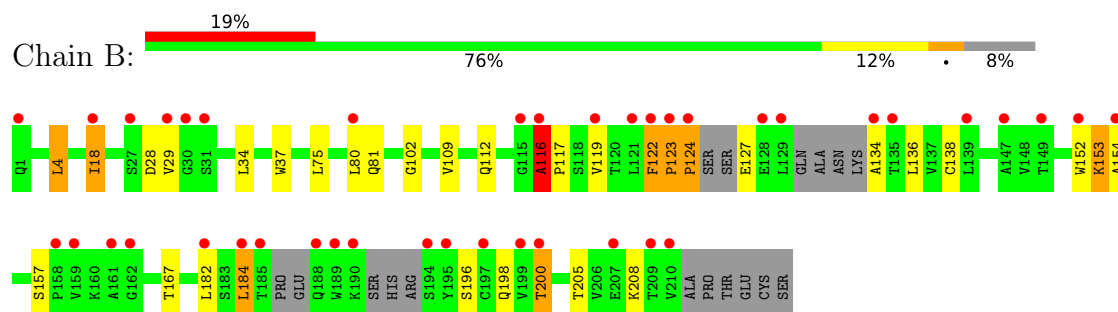
- Molecule 1: DH270.UCA3 heavy chain



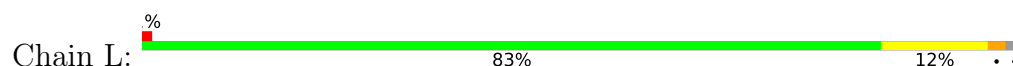
- Molecule 1: DH270.UCA3 heavy chain

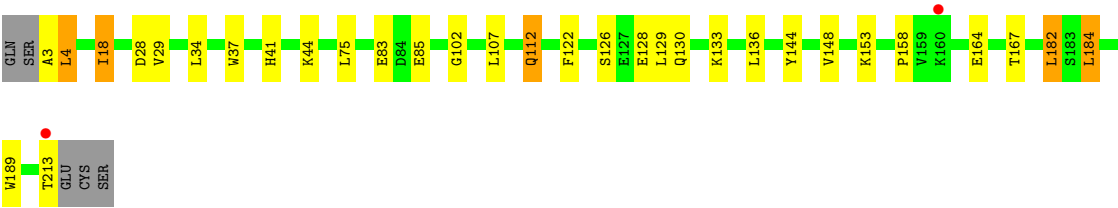


- Molecule 2: DH270.UCA3 light chain



- Molecule 2: DH270.UCA3 light chain





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	98.53Å 76.88Å 137.10Å 90.00° 101.87° 90.00°	Depositor
Resolution (Å)	48.21 – 2.26 48.21 – 2.26	Depositor EDS
% Data completeness (in resolution range)	99.1 (48.21-2.26) 99.1 (48.21-2.26)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 2.27Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.216 , 0.255 0.222 , 0.258	Depositor DCC
R_{free} test set	2355 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	44.0	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.2	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	0.94	EDS
Total number of atoms	6564	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% 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

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	0.39	0/1617	0.84	2/2197 (0.1%)
1	H	0.36	0/1745	0.75	0/2380
2	B	0.42	0/1491	0.90	8/2026 (0.4%)
2	L	0.33	0/1588	0.76	2/2165 (0.1%)
All	All	0.37	0/6441	0.81	12/8768 (0.1%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	104	LEU	N-CA-C	-9.83	97.28	110.55
2	B	124	PRO	CA-N-CD	-8.72	99.79	112.00
2	B	116	ALA	CA-C-N	6.90	143.57	127.00
2	B	116	ALA	C-N-CA	6.90	143.57	127.00
2	B	116	ALA	C-N-CD	-6.50	106.31	120.60
2	L	122	PHE	CA-C-N	5.90	126.45	120.38
2	L	122	PHE	C-N-CA	5.90	126.45	120.38
2	B	122	PHE	CA-C-N	-5.65	114.56	120.38
2	B	122	PHE	C-N-CA	-5.65	114.56	120.38
1	A	223	LYS	N-CA-C	5.19	117.81	108.48
2	B	157	SER	CA-C-N	5.03	124.96	119.78
2	B	157	SER	C-N-CA	5.03	124.96	119.78

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	1579	0	1513	31	0
1	H	1697	0	1639	21	0
2	B	1462	0	1422	28	0
2	L	1552	0	1509	19	0
3	A	70	0	0	2	0
3	B	42	0	0	2	0
3	H	89	0	0	4	0
3	L	73	0	0	3	0
All	All	6564	0	6083	95	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 (95) 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:L:112:GLN:NE2	2:L:144:TYR:OH	1.81	1.14
2:B:122:PHE:C	2:B:124:PRO:HD3	1.80	1.05
1:A:150:ALA:HB2	1:A:200:SER:N	1.84	0.91
1:A:200:SER:O	1:A:202:SER:N	2.07	0.88
2:B:123:PRO:HA	2:B:136:LEU:CD2	2.12	0.79
2:B:123:PRO:HA	2:B:136:LEU:HD23	1.67	0.76
1:H:36:TRP:HD1	1:H:70:MET:HE3	1.50	0.75
2:L:136:LEU:HD12	2:L:182:LEU:HD12	1.68	0.74
2:B:200:THR:HB	2:B:205:THR:HG22	1.69	0.73
1:A:103:SER:OG	1:A:104:LEU:O	2.06	0.73
1:A:173:LEU:HD21	1:A:196:VAL:HG21	1.72	0.72
1:H:36:TRP:CD1	1:H:70:MET:HE3	2.24	0.72
1:A:150:ALA:CB	1:A:200:SER:N	2.54	0.71
1:A:166:VAL:HG12	1:A:212:VAL:HG12	1.75	0.69
2:L:3:ALA:N	3:L:304:HOH:O	2.26	0.68
2:B:153:LYS:HD3	2:B:196:SER:HB2	1.78	0.65
1:A:133:PRO:HB3	1:A:159:TYR:HB3	1.79	0.64
1:A:167:SER:O	1:A:211:ASN:N	2.29	0.63
1:H:213:ASN:HD21	1:H:215:LYS:HG3	1.64	0.62
1:A:135:VAL:C	1:A:136:PHE:HD1	2.08	0.61
2:B:123:PRO:N	2:B:124:PRO:HD3	2.16	0.60
2:B:28:ASP:OD2	2:B:29:VAL:N	2.31	0.60
1:A:206:GLN:HG2	1:A:208:TYR:CZ	2.37	0.59
1:H:200:SER:HA	1:H:203:LEU:HD13	1.83	0.59
2:L:167:THR:OG1	3:L:301:HOH:O	2.17	0.58
2:B:119:VAL:O	2:B:208:LYS:NZ	2.28	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:3:ALA:N	3:L:306:HOH:O	2.37	0.58
2:B:154:ALA:O	3:B:301:HOH:O	2.17	0.57
2:B:122:PHE:O	2:B:124:PRO:HD3	2.03	0.57
1:A:202:SER:OG	1:A:206:GLN:HB2	2.05	0.57
2:L:112:GLN:NE2	2:L:144:TYR:CZ	2.72	0.55
1:A:136:PHE:O	1:A:155:LEU:N	2.29	0.54
2:B:153:LYS:HD2	2:B:198:GLN:CD	2.33	0.53
2:B:122:PHE:HB3	2:B:124:PRO:HD3	1.91	0.53
1:A:166:VAL:HG21	1:A:194:SER:OG	2.08	0.53
1:H:158:ASP:HB3	1:H:189:LEU:HD13	1.91	0.52
1:H:133:PRO:HB3	1:H:159:TYR:HB3	1.93	0.51
1:H:202:SER:OG	3:H:301:HOH:O	2.20	0.51
1:A:177:VAL:HG22	1:A:196:VAL:HB	1.92	0.50
2:L:37:TRP:CD2	2:L:75:LEU:HB2	2.45	0.50
2:L:28:ASP:OD1	2:L:29:VAL:N	2.32	0.49
2:B:18:ILE:HD13	2:B:80:LEU:HD11	1.94	0.49
1:H:184:LEU:HD13	1:H:190:TYR:CE1	2.49	0.48
2:L:4:LEU:HB2	2:L:102:GLY:HA2	1.95	0.48
1:H:101[B]:TRP:HD1	1:H:102:ILE:N	2.12	0.48
1:A:168:TRP:NE1	3:A:309:HOH:O	2.46	0.48
1:A:212:VAL:HG23	1:A:221:VAL:HG23	1.95	0.48
2:B:200:THR:HA	2:B:205:THR:HA	1.95	0.48
1:H:67:ARG:NH1	1:H:90:ASP:OD2	2.34	0.48
1:A:135:VAL:HG21	1:A:212:VAL:HG21	1.96	0.47
2:B:122:PHE:C	2:B:124:PRO:CD	2.71	0.47
1:H:10:GLU:HB2	1:H:123:VAL:HB	1.97	0.47
1:A:40:ALA:HB3	1:A:43:GLN:HG3	1.96	0.46
1:H:157:LYS:NZ	3:H:317:HOH:O	2.48	0.46
1:A:42:GLY:HA3	2:B:167:THR:HG21	1.96	0.46
1:A:138:LEU:HD23	2:B:124:PRO:CG	2.47	0.45
1:A:137:PRO:HD3	1:A:223:LYS:HE2	1.98	0.45
1:H:185:GLN:HG2	2:L:164:GLU:HG3	1.99	0.45
1:H:155:LEU:HD11	1:H:157:LYS:HD3	1.99	0.45
2:L:18:ILE:HD12	2:L:18:ILE:HA	1.79	0.44
2:L:136:LEU:HB2	2:L:182:LEU:HB3	1.99	0.44
2:B:138:CYS:HB2	2:B:152:TRP:CH2	2.53	0.43
1:A:159:TYR:HD1	3:A:301:HOH:O	2.01	0.43
1:H:223:LYS:HA	1:H:223:LYS:HD2	1.79	0.43
2:L:153:LYS:HD3	2:L:158:PRO:HA	2.01	0.43
2:B:134:ALA:O	2:B:184:LEU:N	2.37	0.43
1:H:53:PRO:O	1:H:72:ARG:NE	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:177:VAL:HG22	1:H:196:VAL:HG22	2.00	0.43
1:A:104:LEU:O	1:A:105:TYR:HB2	2.19	0.43
1:A:153:GLY:HA3	1:A:195:VAL:HG12	2.01	0.43
1:A:200:SER:C	1:A:202:SER:N	2.76	0.42
1:A:10:GLU:HB2	1:A:123:VAL:HB	2.00	0.42
2:B:112:GLN:NE2	3:B:306:HOH:O	2.35	0.42
2:B:81:GLN:O	2:B:109:VAL:HG21	2.20	0.42
1:H:36:TRP:CE2	1:H:81:MET:HB2	2.54	0.42
2:L:128:GLU:HG2	2:L:133:LYS:O	2.19	0.42
1:A:206:GLN:HG2	1:A:208:TYR:CE2	2.54	0.42
2:B:4:LEU:HB2	2:B:102:GLY:HA2	2.02	0.42
2:L:126:SER:O	2:L:130:GLN:HG3	2.20	0.42
1:A:152:LEU:HG	1:A:153:GLY:H	1.85	0.41
2:B:123:PRO:N	2:B:124:PRO:CD	2.81	0.41
2:L:83:GLU:H	2:L:83:GLU:CD	2.27	0.41
2:B:116:ALA:HB1	2:B:117:PRO:CA	2.50	0.41
1:H:173:LEU:HD21	1:H:196:VAL:HG11	2.03	0.41
1:A:133:PRO:HB2	1:A:156:VAL:HG23	2.02	0.41
2:B:136:LEU:HB2	2:B:182:LEU:HB3	2.03	0.41
2:L:184:LEU:HD13	2:L:189:TRP:HB2	2.03	0.41
2:L:41:HIS:HB2	2:L:44:LYS:HE2	2.03	0.41
1:A:63:LYS:HE3	1:A:63:LYS:HB2	1.83	0.41
2:B:116:ALA:HB1	2:B:117:PRO:C	2.45	0.41
1:A:138:LEU:HD23	2:B:124:PRO:HG3	2.02	0.40
2:B:37:TRP:CD2	2:B:75:LEU:HB2	2.56	0.40
1:H:185:GLN:NE2	3:H:313:HOH:O	2.53	0.40
2:L:85:GLU:HG3	2:L:107:LEU:O	2.20	0.40
1:H:15:GLY:O	3:H:302:HOH:O	2.22	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	197/238 (83%)	186 (94%)	10 (5%)	1 (0%)	24	24
1	H	218/238 (92%)	211 (97%)	7 (3%)	0	100	100
2	B	189/216 (88%)	180 (95%)	8 (4%)	1 (0%)	24	24
2	L	209/216 (97%)	201 (96%)	8 (4%)	0	100	100
All	All	813/908 (90%)	778 (96%)	33 (4%)	2 (0%)	43	50

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	201	SER
2	B	116	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	172/200 (86%)	163 (95%)	9 (5%)	21	22
1	H	186/200 (93%)	181 (97%)	5 (3%)	39	49
2	B	166/181 (92%)	158 (95%)	8 (5%)	23	25
2	L	176/181 (97%)	167 (95%)	9 (5%)	21	23
All	All	700/762 (92%)	669 (96%)	31 (4%)	25	29

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	104	LEU
1	A	123	VAL
1	A	155	LEU
1	A	156	VAL
1	A	164	VAL
1	A	196	VAL
1	A	212	VAL
1	A	221	VAL
1	A	224	ARG

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Mol	Chain	Res	Type
2	B	4	LEU
2	B	18	ILE
2	B	34	LEU
2	B	123	PRO
2	B	127	GLU
2	B	153	LYS
2	B	184	LEU
2	B	200	THR
1	H	1	GLN
1	H	123	VAL
1	H	155	LEU
1	H	205	THR
1	H	221	VAL
2	L	4	LEU
2	L	18	ILE
2	L	34	LEU
2	L	112	GLN
2	L	129	LEU
2	L	148	VAL
2	L	182	LEU
2	L	184	LEU
2	L	213	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	54	ASN
1	A	113	ASN
2	B	81	GLN
2	B	112	GLN
2	B	171	GLN
2	B	188	GLN
1	H	43	GLN
1	H	213	ASN
2	L	112	GLN
2	L	198	GLN

5.3.3 RNA ⓘ

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

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	207/238 (86%)	1.23	57 (27%) 1 1	28, 57, 140, 185	0
1	H	221/238 (92%)	0.36	8 (3%) 46 46	22, 45, 71, 100	1 (0%)
2	B	199/216 (92%)	1.11	41 (20%) 2 2	30, 57, 124, 188	0
2	L	211/216 (97%)	0.20	2 (0%) 81 82	26, 44, 66, 105	0
All	All	838/908 (92%)	0.71	108 (12%) 7 7	22, 48, 120, 188	1 (0%)

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	152	LEU	7.9
2	B	124	PRO	7.5
1	A	137	PRO	5.9
1	A	155	LEU	5.8
2	B	134	ALA	5.6
2	B	185	THR	5.6
1	A	135	VAL	5.6
1	A	151	ALA	5.4
1	A	209	ILE	5.4
1	A	214	HIS	5.3
1	A	168	TRP	5.0
1	A	208	TYR	4.9
1	A	134	SER	4.8
2	B	115	GLY	4.8
1	A	176	GLY	4.8
1	A	154	CYS	4.7
2	B	190	LYS	4.7
1	A	202	SER	4.6
1	A	136	PHE	4.6
2	B	129	LEU	4.6
1	A	172	ALA	4.5

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Mol	Chain	Res	Type	RSRZ
2	B	122	PHE	4.5
1	A	221	VAL	4.5
1	A	200	SER	4.5
1	A	210	CYS	4.4
1	A	195	VAL	4.4
1	A	213	ASN	4.2
1	A	196	VAL	4.1
1	A	222	ASP	4.1
1	A	173	LEU	4.0
2	B	116	ALA	4.0
1	A	156	VAL	4.0
2	B	159	VAL	4.0
1	A	166	VAL	3.9
1	A	207	THR	3.9
2	B	195	TYR	3.9
2	B	184	LEU	3.8
2	B	189	TRP	3.6
1	A	205	THR	3.6
1	A	175	SER	3.6
2	B	194	SER	3.5
1	A	138	LEU	3.5
2	B	135	THR	3.4
2	B	209	THR	3.3
1	A	194	SER	3.3
1	A	129	SER	3.2
2	B	210	VAL	3.2
1	A	161	PRO	3.2
1	A	149	THR	3.2
2	B	121	LEU	3.1
2	B	161	ALA	3.1
2	B	123	PRO	3.1
1	A	174	THR	3.1
1	A	150	ALA	3.1
1	A	169	ASN	3.0
1	H	228	LYS	3.0
1	H	141	SER	3.0
2	B	31	SER	3.0
2	B	1	GLN	2.9
1	A	201	SER	2.9
1	H	149	THR	2.8
1	A	212	VAL	2.8
2	B	30	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	163	PRO	2.7
2	B	29	VAL	2.7
2	B	18	ILE	2.7
1	A	223	LYS	2.6
2	B	128	GLU	2.6
2	B	162	GLY	2.6
2	B	197	CYS	2.6
1	A	159	TYR	2.6
1	A	217	SER	2.5
1	A	177	VAL	2.5
1	A	165	THR	2.4
2	B	139	LEU	2.4
2	B	149	THR	2.4
1	A	164	VAL	2.4
1	A	153	GLY	2.4
1	H	172	ALA	2.4
2	B	207	GLU	2.4
1	A	178	HIS	2.3
1	A	171	GLY	2.3
1	A	224	ARG	2.3
2	B	158	PRO	2.3
2	B	154	ALA	2.3
2	B	182	LEU	2.3
2	B	152	TRP	2.3
1	A	62	GLN	2.2
2	B	199	VAL	2.2
1	A	133	PRO	2.2
2	B	80	LEU	2.2
2	B	119	VAL	2.2
2	B	27	SER	2.2
2	B	147	ALA	2.2
2	B	200	THR	2.2
1	A	130	THR	2.1
1	A	170	SER	2.1
1	H	17	SER	2.1
1	A	59	ASN	2.1
1	H	54	ASN	2.1
2	B	188	GLN	2.1
1	A	188	GLY	2.1
1	H	66	GLY	2.1
1	A	1	GLN	2.1
2	L	213	THR	2.0

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Mol	Chain	Res	Type	RSRZ
2	L	160	LYS	2.0
1	H	62	GLN	2.0
1	A	211	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.