



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:55 AM UTC

PDB ID : 6XV8 / pdb_00006xv8
Title : Crystal structure of Megabody Mb-Nb207-c7HopQ_G10
Authors : Steyaert, J.; Uchanski, T.; Fischer, B.
Deposited on : 2020-01-21
Resolution : 3.15 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

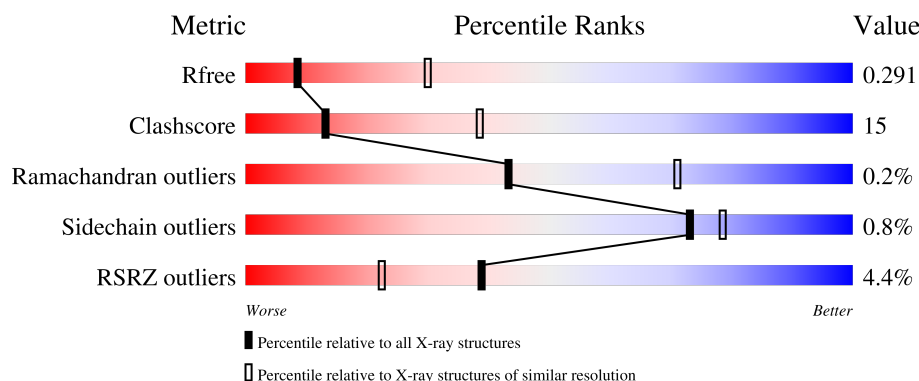
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 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	2361 (3.20-3.12)
Clashscore	190562	2486 (3.20-3.12)
Ramachandran outliers	187476	2405 (3.20-3.12)
Sidechain outliers	187428	2404 (3.20-3.12)
RSRZ outliers	180081	2361 (3.20-3.12)

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	521	3% 58%21%21%
1	B	521	3% 57%25%18%
1	C	521	3% 58%21%21%
1	D	521	5% 51%22%26%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 12431 atoms, of which 0 are hydrogens and 0 are deuteriums.

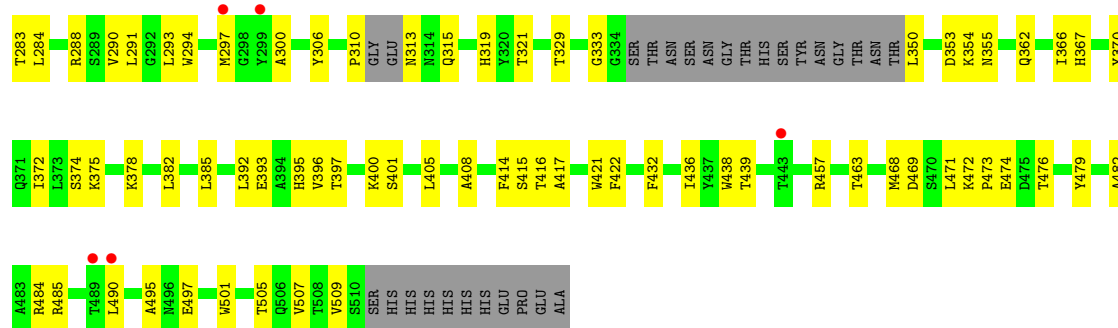
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Outer membrane protein.

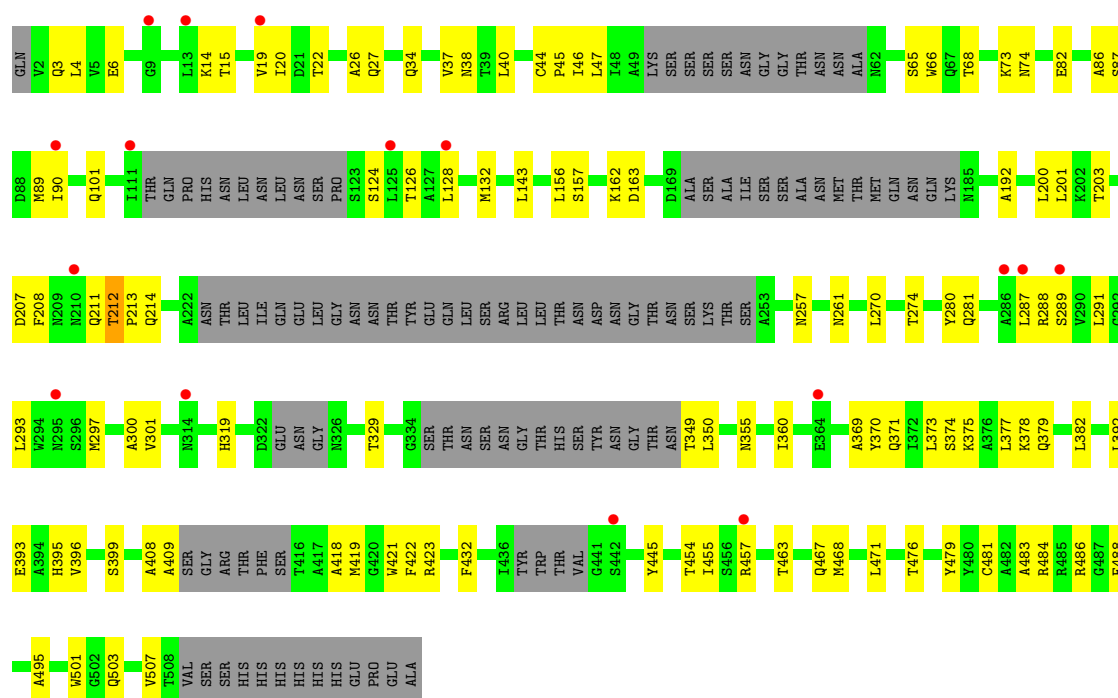
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	412	Total	C	N	O	S	0	0	0
			3134	1942	552	627	13			
1	B	429	Total	C	N	O	S	0	0	0
			3251	2012	572	654	13			
1	C	412	Total	C	N	O	S	0	0	0
			3112	1924	549	626	13			
1	D	387	Total	C	N	O	S	0	0	0
			2934	1820	514	588	12			

There are 8 discrepancies between the modelled and reference sequences:

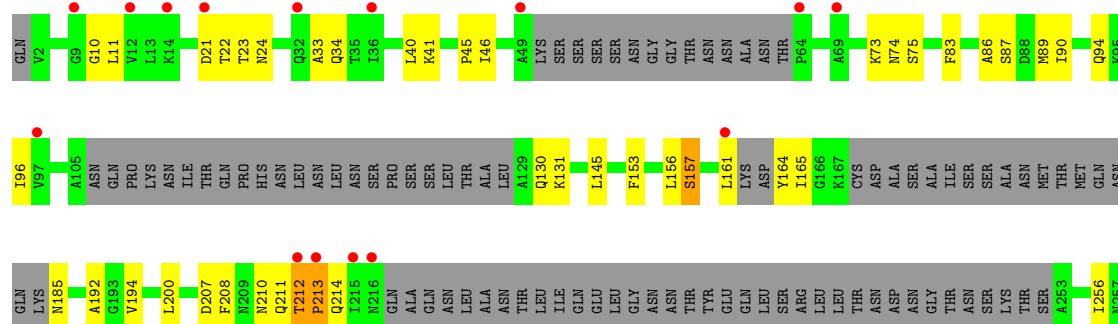
Chain	Residue	Modelled	Actual	Comment	Reference
A	13	LEU	-	linker	UNP B5Z8H1
A	401	SER	-	linker	UNP B5Z8H1
B	13	LEU	-	linker	UNP B5Z8H1
B	401	SER	-	linker	UNP B5Z8H1
C	13	LEU	-	linker	UNP B5Z8H1
C	401	SER	-	linker	UNP B5Z8H1
D	13	LEU	-	linker	UNP B5Z8H1
D	401	SER	-	linker	UNP B5Z8H1

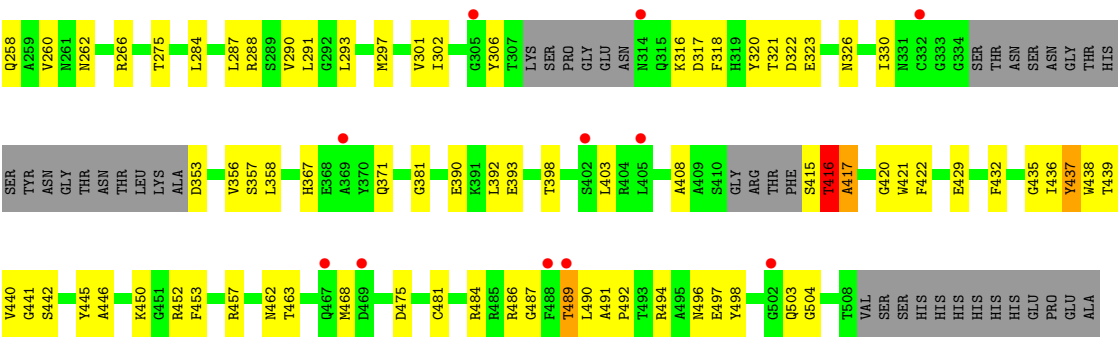


• Molecule 1: Outer membrane protein



• Molecule 1: Outer membrane protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.24Å 155.05Å 89.75Å 90.00° 91.73° 90.00°	Depositor
Resolution (Å)	79.21 – 3.15 79.21 – 3.15	Depositor EDS
% Data completeness (in resolution range)	96.0 (79.21-3.15) 96.1 (79.21-3.15)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 3.13Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.255 , 0.297 0.252 , 0.291	Depositor DCC
R_{free} test set	1686 reflections (4.50%)	wwPDB-VP
Wilson B-factor (Å ²)	101.5	Xtriage
Anisotropy	0.207	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 79.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.001 for h,-k,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	12431	wwPDB-VP
Average B, all atoms (Å ²)	114.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.98% 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	0.43	0/3178	0.74	1/4298 (0.0%)
1	B	0.49	0/3296	0.86	4/4460 (0.1%)
1	C	0.52	0/3151	0.84	3/4258 (0.1%)
1	D	0.42	0/2974	0.84	8/4019 (0.2%)
All	All	0.47	0/12599	0.82	16/17035 (0.1%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	107	GLN	CA-C-N	12.29	132.69	120.52
1	B	107	GLN	C-N-CA	12.29	132.69	120.52
1	D	212	THR	CA-C-N	7.84	129.64	119.84
1	D	212	THR	C-N-CA	7.84	129.64	119.84
1	D	417	ALA	N-CA-C	7.63	119.24	111.07
1	C	457	ARG	N-CA-C	7.53	121.87	109.06
1	B	164	TYR	N-CA-C	7.39	122.68	112.04
1	D	11	LEU	N-CA-C	-6.48	98.17	108.73
1	B	70	GLY	N-CA-C	6.46	120.48	112.73
1	D	212	THR	O-C-N	6.26	124.41	121.53
1	D	416	THR	N-CA-C	5.73	119.25	112.72
1	C	212	THR	CA-C-N	5.38	126.56	119.84
1	C	212	THR	C-N-CA	5.38	126.56	119.84
1	D	157	SER	N-CA-C	5.21	119.38	113.19
1	A	475	ASP	N-CA-C	-5.13	99.88	110.80
1	D	489	THR	N-CA-C	5.13	116.95	111.36

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	3134	0	3069	82	0
1	B	3251	0	3177	116	0
1	C	3112	0	3046	82	0
1	D	2934	0	2856	97	0
All	All	12431	0	12148	366	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 15.

All (366) 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:B:211:GLN:HG3	1:B:266:ARG:NH2	1.68	1.07
1:D:256:ILE:O	1:D:260:VAL:HG23	1.62	0.98
1:B:108:PRO:HD2	1:B:264:ASN:OD1	1.75	0.87
1:B:7:SER:HB2	1:C:101:GLN:HE22	1.44	0.82
1:A:14:LYS:HG2	1:A:400:LYS:HB2	1.62	0.82
1:B:468:MET:HE2	1:B:471:LEU:HD21	1.62	0.80
1:B:21:ASP:OD1	1:B:393:GLU:HG2	1.82	0.79
1:D:46:ILE:HA	1:D:74:ASN:HA	1.64	0.79
1:B:34:GLN:HG3	1:B:90:ILE:HG21	1.64	0.77
1:A:479:TYR:HE2	1:A:507:VAL:HG11	1.51	0.76
1:D:353:ASP:HB2	1:D:357:SER:O	1.85	0.76
1:D:266:ARG:HA	1:D:266:ARG:NE	2.00	0.76
1:B:211:GLN:HG3	1:B:266:ARG:CZ	2.15	0.75
1:D:22:THR:O	1:D:392:LEU:HD23	1.87	0.74
1:A:476:THR:HG23	1:A:508:THR:HA	1.71	0.73
1:B:161:LEU:O	1:B:165:ILE:HG22	1.90	0.72
1:B:422:PHE:HE2	1:B:495:ALA:HA	1.53	0.72
1:C:74:ASN:ND2	1:C:355:ASN:OD1	2.23	0.71
1:C:214:GLN:N	1:C:214:GLN:OE1	2.21	0.71
1:C:468:MET:HE2	1:C:471:LEU:HD21	1.70	0.71
1:D:192:ALA:HB1	1:D:288:ARG:HB2	1.72	0.70
1:C:86:ALA:HA	1:C:89:MET:HE3	1.72	0.70
1:D:185:ASN:N	1:D:326:ASN:HD22	1.89	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:6:GLU:CD	1:C:6:GLU:H	2.00	0.69
1:B:422:PHE:CE1	1:B:432:PHE:HD1	2.11	0.68
1:B:306:TYR:CD1	1:B:350:LEU:O	2.46	0.68
1:B:306:TYR:HD1	1:B:350:LEU:O	1.77	0.67
1:B:14:LYS:HB3	1:C:20:ILE:HG12	1.75	0.67
1:C:46:ILE:HA	1:C:74:ASN:HA	1.76	0.67
1:A:375:LYS:HE3	1:A:379:GLN:OE1	1.94	0.67
1:B:422:PHE:HE1	1:B:432:PHE:HD1	1.41	0.67
1:D:484:ARG:HE	1:D:486:ARG:HB2	1.60	0.67
1:B:12:VAL:HA	1:C:22:THR:HG22	1.76	0.67
1:D:207:ASP:O	1:D:211:GLN:NE2	2.29	0.66
1:D:415:SER:HB3	1:D:462:ASN:ND2	2.11	0.66
1:D:438:TRP:O	1:D:441:GLY:N	2.26	0.65
1:D:322:ASP:HB3	1:D:326:ASN:H	1.62	0.65
1:B:416:THR:HG22	1:B:439:THR:HG22	1.78	0.65
1:D:214:GLN:HG3	1:D:214:GLN:O	1.97	0.65
1:B:7:SER:HB2	1:C:101:GLN:NE2	2.11	0.65
1:C:6:GLU:OE2	1:C:503:GLN:N	2.30	0.65
1:D:416:THR:O	1:D:438:TRP:HA	1.96	0.64
1:A:414:PHE:CD2	1:A:483:ALA:HB1	2.32	0.64
1:B:438:TRP:O	1:B:457:ARG:NH1	2.31	0.64
1:D:10:GLY:O	1:D:403:LEU:HD23	1.98	0.64
1:B:283:THR:CG2	1:B:382:LEU:HD21	2.28	0.63
1:B:7:SER:HB3	1:C:101:GLN:OE1	1.98	0.63
1:B:422:PHE:HE1	1:B:432:PHE:CD1	2.16	0.63
1:D:258:GLN:HA	1:D:258:GLN:OE1	1.98	0.63
1:A:476:THR:HA	1:A:507:VAL:HG13	1.80	0.63
1:D:408:ALA:HA	1:D:463:THR:HG22	1.81	0.62
1:B:185:ASN:OD1	1:B:186:ASN:N	2.32	0.62
1:C:484:ARG:HE	1:C:486:ARG:HB2	1.65	0.62
1:A:474:GLU:CD	1:A:474:GLU:O	2.43	0.62
1:C:257:ASN:O	1:C:261:ASN:ND2	2.33	0.61
1:C:128:LEU:HG	1:C:132:MET:HE2	1.81	0.61
1:A:315:GLN:HG2	1:A:469:ASP:HB3	1.83	0.61
1:A:479:TYR:CE2	1:A:507:VAL:HG11	2.35	0.61
1:C:192:ALA:HB1	1:C:288:ARG:HB2	1.83	0.61
1:A:477:ALA:O	1:A:507:VAL:HG12	2.01	0.60
1:C:291:LEU:HD12	1:C:373:LEU:HG	1.83	0.60
1:B:473:PRO:HA	1:B:509:VAL:HB	1.84	0.60
1:D:322:ASP:HB3	1:D:326:ASN:N	2.16	0.60
1:A:108:PRO:HB2	1:A:260:VAL:HG22	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:304:GLY:O	1:A:352:ALA:HB2	2.03	0.59
1:C:89:MET:HE1	1:C:156:LEU:HD22	1.83	0.59
1:A:34:GLN:HG2	1:A:38:ASN:ND2	2.17	0.59
1:B:315:GLN:HB3	1:B:333:GLY:HA2	1.85	0.58
1:B:90:ILE:HD11	1:B:290:VAL:HG21	1.85	0.58
1:A:438:TRP:O	1:A:457:ARG:NH2	2.34	0.58
1:B:34:GLN:HG3	1:B:90:ILE:CG2	2.34	0.58
1:B:204:SER:HB2	1:B:276:ASN:O	2.04	0.58
1:C:212:THR:N	1:C:213:PRO:HD2	2.19	0.58
1:C:163:ASP:OD1	1:C:163:ASP:N	2.31	0.58
1:B:354:LYS:O	1:B:355:ASN:HB2	2.03	0.58
1:A:109:LYS:HG3	1:A:110:ASN:H	1.69	0.58
1:D:200:LEU:HD13	1:D:275:THR:HG23	1.86	0.58
1:A:185:ASN:OD1	1:A:186:ASN:N	2.36	0.58
1:D:46:ILE:HG23	1:D:73:LYS:O	2.04	0.57
1:D:212:THR:O	1:D:212:THR:HG22	2.04	0.57
1:A:130:GLN:HA	1:A:133:LEU:HB3	1.86	0.57
1:B:476:THR:HG23	1:B:507:VAL:O	2.03	0.57
1:B:291:LEU:HD21	1:B:374:SER:HA	1.86	0.57
1:D:24:ASN:ND2	1:D:381:GLY:O	2.35	0.57
1:B:125:LEU:N	1:B:125:LEU:HD22	2.20	0.57
1:C:495:ALA:O	1:C:501:TRP:HZ2	1.88	0.57
1:B:8:GLY:HA3	1:B:405:LEU:HD23	1.87	0.56
1:A:106:ASN:HD22	1:A:131:LYS:HE2	1.70	0.56
1:C:44:CYS:O	1:C:74:ASN:HB2	2.04	0.56
1:C:291:LEU:HD13	1:C:377:LEU:HD12	1.87	0.56
1:B:192:ALA:HB1	1:B:288:ARG:HG3	1.87	0.56
1:D:21:ASP:OD1	1:D:393:GLU:HG2	2.05	0.56
1:D:492:PRO:HB2	1:D:498:TYR:CE2	2.41	0.56
1:A:476:THR:HG22	1:A:476:THR:O	2.04	0.56
1:B:200:LEU:HA	1:B:203:THR:HG22	1.88	0.56
1:A:108:PRO:HD3	1:A:135:ASN:HD22	1.71	0.56
1:C:375:LYS:HE2	1:C:379:GLN:OE1	2.06	0.56
1:D:452:ARG:HH11	1:D:475:ASP:CG	2.14	0.56
1:C:26:ALA:HB2	1:C:382:LEU:HD11	1.88	0.56
1:D:266:ARG:HA	1:D:266:ARG:HE	1.70	0.56
1:A:22:THR:O	1:A:392:LEU:HG	2.06	0.55
1:A:421:TRP:CH2	1:A:481:CYS:HB3	2.40	0.55
1:B:125:LEU:HD22	1:B:125:LEU:H	1.71	0.55
1:A:453:PHE:CE1	1:A:468:MET:HE3	2.42	0.55
1:D:96:ILE:HD11	1:D:145:LEU:HB2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:318:PHE:HD1	1:D:367:HIS:ND1	2.04	0.55
1:B:210:ASN:O	1:B:213:PRO:HD2	2.07	0.55
1:B:9:GLY:HA3	1:C:27:GLN:CG	2.37	0.55
1:B:212:THR:O	1:B:216:ASN:HB2	2.07	0.54
1:D:306:TYR:C	1:D:306:TYR:CD2	2.85	0.54
1:D:484:ARG:HD2	1:D:497:GLU:O	2.06	0.54
1:D:432:PHE:O	1:D:446:ALA:HB2	2.08	0.54
1:B:414:PHE:O	1:B:417:ALA:N	2.32	0.54
1:C:4:LEU:HB2	1:C:501:TRP:O	2.08	0.54
1:D:208:PHE:CE1	1:D:266:ARG:HD3	2.42	0.54
1:B:161:LEU:HB3	1:B:165:ILE:CG2	2.38	0.54
1:B:143:LEU:HD11	1:B:208:PHE:HB3	1.90	0.53
1:D:23:THR:HG22	1:D:390:GLU:C	2.33	0.53
1:B:306:TYR:C	1:B:306:TYR:CD2	2.86	0.53
1:B:19:VAL:O	1:C:14:LYS:HA	2.08	0.53
1:D:284:LEU:O	1:D:288:ARG:HG2	2.08	0.53
1:D:441:GLY:CA	1:D:457:ARG:HH12	2.21	0.53
1:D:320:TYR:HE2	1:D:330:ILE:HD12	1.74	0.53
1:A:130:GLN:O	1:A:134:LYS:N	2.24	0.53
1:C:208:PHE:CE2	1:C:270:LEU:HD13	2.43	0.53
1:D:208:PHE:HE1	1:D:266:ARG:HD3	1.73	0.53
1:A:452:ARG:HG3	1:A:453:PHE:HD2	1.73	0.53
1:B:43:TYR:HE2	1:B:362:GLN:HG3	1.73	0.53
1:A:192:ALA:HB3	1:A:194:VAL:HG13	1.91	0.53
1:A:422:PHE:HE2	1:A:498:TYR:CD2	2.26	0.53
1:A:453:PHE:CE1	1:A:468:MET:HG2	2.43	0.53
1:A:457:ARG:O	1:A:457:ARG:HG2	2.09	0.53
1:B:414:PHE:O	1:B:415:SER:C	2.51	0.53
1:A:215:ILE:HA	1:A:218:ALA:HB3	1.90	0.53
1:C:289:SER:O	1:C:293:LEU:N	2.37	0.52
1:D:34:GLN:NE2	1:D:94:GLN:OE1	2.42	0.52
1:B:315:GLN:O	1:B:400:LYS:HE2	2.09	0.52
1:D:452:ARG:NH1	1:D:475:ASP:OD1	2.42	0.52
1:B:19:VAL:HG22	1:B:395:HIS:CD2	2.44	0.52
1:D:494:ARG:HD2	1:D:496:ASN:OD1	2.08	0.52
1:C:349:THR:OG1	1:C:350:LEU:N	2.40	0.52
1:C:371:GLN:HB2	1:C:396:VAL:HG22	1.90	0.52
1:D:322:ASP:OD1	1:D:323:GLU:N	2.42	0.52
1:A:368:GLU:HG3	1:A:396:VAL:HG21	1.93	0.51
1:D:422:PHE:CD1	1:D:432:PHE:HA	2.44	0.51
1:A:316:LYS:HD2	1:A:398:THR:HB	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:422:PHE:CE2	1:B:495:ALA:HA	2.41	0.51
1:D:33:ALA:HB1	1:D:290:VAL:HG21	1.91	0.51
1:A:46:ILE:HA	1:A:74:ASN:HA	1.93	0.51
1:B:297:MET:SD	1:B:297:MET:C	2.94	0.51
1:D:437:TYR:CZ	1:D:491:ALA:HB1	2.45	0.51
1:D:489:THR:HG22	1:D:490:LEU:HG	1.92	0.51
1:D:415:SER:HB3	1:D:462:ASN:CG	2.36	0.51
1:A:135:ASN:O	1:A:139:GLN:HG3	2.10	0.51
1:B:22:THR:O	1:B:392:LEU:HD23	2.12	0.50
1:A:103:LEU:HD11	1:A:267:ALA:HB1	1.93	0.50
1:A:212:THR:HB	1:A:213:PRO:HD3	1.92	0.50
1:A:317:ASP:HB2	1:A:399:SER:HB3	1.93	0.50
1:B:44:CYS:O	1:B:74:ASN:HB2	2.11	0.50
1:C:46:ILE:HG23	1:C:73:LYS:O	2.12	0.50
1:C:445:TYR:OH	1:C:455:ILE:HG22	2.12	0.50
1:D:317:ASP:HA	1:D:330:ILE:O	2.12	0.50
1:C:476:THR:HG23	1:C:507:VAL:O	2.12	0.50
1:B:143:LEU:CD1	1:B:208:PHE:HB3	2.42	0.50
1:B:315:GLN:NE2	1:B:469:ASP:HB2	2.27	0.50
1:A:5:VAL:HG23	1:A:408:ALA:HB3	1.93	0.50
1:A:18:SER:HB2	1:A:396:VAL:HB	1.94	0.50
1:C:22:THR:O	1:C:392:LEU:HD23	2.12	0.50
1:B:46:ILE:HA	1:B:74:ASN:HA	1.94	0.49
1:B:294:TRP:CH2	1:B:366:ILE:HG21	2.47	0.49
1:B:422:PHE:CE1	1:B:432:PHE:CD1	2.95	0.49
1:C:207:ASP:O	1:C:211:GLN:HG2	2.12	0.49
1:B:408:ALA:HA	1:B:463:THR:HG22	1.94	0.49
1:A:14:LYS:HE2	1:A:16:THR:CG2	2.42	0.49
1:A:417:ALA:HB1	1:A:484:ARG:O	2.12	0.49
1:C:200:LEU:HA	1:C:203:THR:HG22	1.94	0.49
1:B:421:TRP:O	1:B:422:PHE:HD1	1.95	0.49
1:C:374:SER:O	1:C:378:LYS:HG2	2.11	0.49
1:A:261:ASN:O	1:A:265:GLU:HB2	2.12	0.49
1:B:415:SER:O	1:B:457:ARG:NH1	2.44	0.49
1:B:47:LEU:HA	1:B:300:ALA:O	2.13	0.48
1:B:208:PHE:CE2	1:B:270:LEU:HD13	2.48	0.48
1:B:22:THR:HB	1:B:28:ASN:ND2	2.28	0.48
1:C:40:LEU:CD1	1:C:297:MET:HE1	2.43	0.48
1:B:283:THR:HG21	1:B:385:LEU:HD21	1.94	0.48
1:C:47:LEU:HA	1:C:300:ALA:O	2.13	0.48
1:C:319:HIS:ND1	1:C:329:THR:HG22	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:ILE:HG13	1:B:112:THR:H	1.79	0.48
1:B:479:TYR:O	1:B:505:THR:HG22	2.13	0.48
1:D:452:ARG:HG3	1:D:453:PHE:CD1	2.49	0.48
1:A:291:LEU:HD12	1:A:370:TYR:CE1	2.49	0.48
1:D:157:SER:OG	1:D:165:ILE:HG21	2.14	0.48
1:A:422:PHE:CD1	1:A:432:PHE:HA	2.49	0.48
1:C:408:ALA:HA	1:C:463:THR:HG22	1.96	0.48
1:B:150:GLU:HB2	1:B:201:LEU:CD2	2.44	0.48
1:C:143:LEU:HD23	1:C:143:LEU:HA	1.69	0.47
1:D:161:LEU:HD12	1:D:164:TYR:HB2	1.95	0.47
1:D:320:TYR:CE2	1:D:330:ILE:HD12	2.49	0.47
1:D:153:PHE:O	1:D:156:LEU:HB3	2.14	0.47
1:A:156:LEU:HD11	1:A:293:LEU:HD12	1.95	0.47
1:B:12:VAL:O	1:B:12:VAL:HG23	2.13	0.47
1:C:201:LEU:HD13	1:C:281:GLN:HB2	1.97	0.47
1:C:422:PHE:CD1	1:C:432:PHE:HA	2.49	0.47
1:A:452:ARG:HG3	1:A:453:PHE:CD2	2.50	0.47
1:D:130:GLN:HG2	1:D:131:LYS:N	2.30	0.47
1:B:9:GLY:HA3	1:C:27:GLN:HG3	1.97	0.47
1:B:39:THR:OG1	1:B:366:ILE:HD11	2.15	0.47
1:C:6:GLU:N	1:C:6:GLU:OE1	2.47	0.47
1:C:212:THR:HB	1:C:213:PRO:HD3	1.96	0.47
1:C:454:THR:HB	1:C:467:GLN:HB3	1.97	0.47
1:D:287:LEU:O	1:D:290:VAL:HG22	2.14	0.47
1:D:417:ALA:HB3	1:D:484:ARG:O	2.15	0.47
1:D:484:ARG:NH2	1:D:487:GLY:O	2.47	0.47
1:B:161:LEU:HB3	1:B:165:ILE:HG22	1.97	0.47
1:B:319:HIS:ND1	1:B:329:THR:HG22	2.29	0.47
1:D:302:ILE:O	1:D:357:SER:HB2	2.15	0.47
1:D:422:PHE:CE1	1:D:432:PHE:HB2	2.50	0.47
1:D:437:TYR:CE2	1:D:491:ALA:HB1	2.50	0.47
1:B:111:ILE:HG13	1:B:112:THR:N	2.30	0.47
1:A:149:VAL:O	1:A:153:PHE:N	2.43	0.47
1:B:150:GLU:HB2	1:B:201:LEU:HD23	1.97	0.47
1:A:436:ILE:HA	1:A:442:SER:O	2.15	0.47
1:B:36:ILE:HD11	1:B:370:TYR:HA	1.97	0.47
1:D:445:TYR:HB2	1:D:450:LYS:HE2	1.97	0.47
1:A:422:PHE:CE1	1:A:432:PHE:HD1	2.34	0.46
1:C:15:THR:HG22	1:C:399:SER:HB2	1.96	0.46
1:D:161:LEU:HA	1:D:164:TYR:HD2	1.80	0.46
1:A:379:GLN:HG2	1:B:438:TRP:CD1	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:306:TYR:CZ	1:B:310:PRO:HB3	2.50	0.46
1:C:65:SER:O	1:C:68:THR:HG22	2.15	0.46
1:D:429:GLU:HA	1:D:429:GLU:OE1	2.15	0.46
1:D:161:LEU:HD13	1:D:164:TYR:CD2	2.50	0.46
1:A:414:PHE:N	1:A:414:PHE:CD1	2.84	0.46
1:A:449:ALA:HB1	1:A:453:PHE:HB2	1.97	0.46
1:B:472:LYS:HB3	1:B:474:GLU:OE1	2.16	0.46
1:D:421:TRP:CZ3	1:D:481:CYS:HB3	2.49	0.46
1:B:319:HIS:CD2	1:B:397:THR:HG22	2.51	0.46
1:B:212:THR:HB	1:B:213:PRO:HD3	1.96	0.46
1:D:439:THR:HA	1:D:457:ARG:NH2	2.31	0.46
1:D:420:GLY:HA2	1:D:435:GLY:HA2	1.97	0.46
1:C:408:ALA:HB2	1:C:463:THR:HG22	1.97	0.46
1:A:79:PHE:HA	1:A:82:GLU:OE1	2.16	0.45
1:A:142:ILE:HG21	1:A:270:LEU:HD21	1.98	0.45
1:A:421:TRP:CZ3	1:A:481:CYS:HB3	2.51	0.45
1:B:375:LYS:HD3	1:B:378:LYS:HE2	1.99	0.45
1:A:11:LEU:C	1:A:11:LEU:HD23	2.41	0.45
1:B:482:ALA:HB2	1:B:501:TRP:CE3	2.51	0.45
1:D:316:LYS:HD3	1:D:398:THR:HB	1.98	0.45
1:A:89:MET:HE1	1:A:156:LEU:HD22	1.99	0.45
1:C:157:SER:O	1:C:162:LYS:HA	2.17	0.45
1:C:287:LEU:HB3	1:C:377:LEU:HD11	1.97	0.45
1:D:192:ALA:HB3	1:D:194:VAL:HG13	1.99	0.45
1:D:403:LEU:HD12	1:D:468:MET:SD	2.57	0.45
1:D:422:PHE:HE1	1:D:432:PHE:HB2	1.81	0.45
1:D:437:TYR:CD1	1:D:437:TYR:N	2.84	0.45
1:B:26:ALA:HB2	1:B:382:LEU:HD11	1.99	0.45
1:B:107:GLN:O	1:B:107:GLN:HG2	2.17	0.45
1:C:37:VAL:HG11	1:C:87:SER:HA	1.98	0.45
1:D:45:PRO:O	1:D:75:SER:HB2	2.16	0.45
1:D:318:PHE:HB3	1:D:367:HIS:CE1	2.52	0.45
1:C:40:LEU:HD13	1:C:297:MET:HE1	1.99	0.45
1:B:313:ASN:N	1:B:313:ASN:OD1	2.50	0.44
1:B:260:VAL:HG13	1:B:264:ASN:ND2	2.33	0.44
1:B:414:PHE:HE1	1:B:485:ARG:N	2.16	0.44
1:B:436:ILE:HD11	1:B:457:ARG:HH21	1.82	0.44
1:D:367:HIS:O	1:D:371:GLN:HG2	2.17	0.44
1:A:475:ASP:OD2	1:A:475:ASP:N	2.50	0.44
1:A:110:ASN:HB2	1:A:260:VAL:HG11	1.98	0.44
1:B:306:TYR:CD2	1:B:306:TYR:O	2.70	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:19:VAL:HG22	1:C:395:HIS:CE1	2.53	0.44
1:C:34:GLN:HG2	1:C:38:ASN:ND2	2.32	0.44
1:C:445:TYR:CZ	1:C:455:ILE:HG22	2.53	0.44
1:D:503:GLN:HG3	1:D:504:GLY:O	2.17	0.44
1:D:34:GLN:HB3	1:D:90:ILE:HG21	2.00	0.44
1:C:3:GLN:O	1:C:409:ALA:HA	2.17	0.44
1:A:47:LEU:HD22	1:A:75:SER:HB3	2.00	0.44
1:A:193:GLY:O	1:A:288:ARG:NH1	2.51	0.44
1:B:257:ASN:O	1:B:261:ASN:ND2	2.51	0.44
1:A:453:PHE:HE1	1:A:468:MET:HE3	1.80	0.43
1:A:468:MET:HB3	1:A:471:LEU:HD21	1.99	0.43
1:D:301:VAL:HG11	1:D:358:LEU:HD12	2.00	0.43
1:B:144:LYS:HE2	1:B:144:LYS:HB3	1.79	0.43
1:D:86:ALA:HA	1:D:89:MET:CE	2.48	0.43
1:B:90:ILE:HD11	1:B:290:VAL:CG2	2.48	0.43
1:C:418:ALA:HB2	1:C:488:PHE:CE1	2.53	0.43
1:D:208:PHE:HE1	1:D:266:ARG:CD	2.30	0.43
1:B:319:HIS:HD1	1:B:329:THR:HG22	1.83	0.43
1:C:274:THR:O	1:C:281:GLN:NE2	2.41	0.43
1:B:43:TYR:CE2	1:B:362:GLN:HG3	2.52	0.43
1:C:408:ALA:CA	1:C:463:THR:HG22	2.49	0.43
1:A:86:ALA:HA	1:A:89:MET:CE	2.49	0.43
1:B:164:TYR:HB3	1:B:188:GLY:O	2.19	0.43
1:D:40:LEU:HD13	1:D:83:PHE:CZ	2.53	0.43
1:C:124:SER:C	1:C:126:THR:H	2.26	0.43
1:A:14:LYS:HD3	1:A:400:LYS:HD2	2.01	0.43
1:A:136:ALA:HB2	1:A:263:LEU:HD21	2.00	0.43
1:C:423:ARG:HD3	1:C:479:TYR:OH	2.18	0.43
1:C:419:MET:HE3	1:C:483:ALA:HB2	2.00	0.42
1:A:492:PRO:HB2	1:A:498:TYR:CE1	2.54	0.42
1:B:108:PRO:HG2	1:B:260:VAL:HG13	2.00	0.42
1:B:211:GLN:CG	1:B:266:ARG:NH2	2.61	0.42
1:B:490:LEU:HD23	1:B:490:LEU:HA	1.85	0.42
1:C:90:ILE:HD13	1:C:90:ILE:HA	1.80	0.42
1:C:280:TYR:OH	1:C:382:LEU:O	2.22	0.42
1:D:41:LYS:NZ	1:D:87:SER:HB2	2.34	0.42
1:A:407:CYS:O	1:A:463:THR:HA	2.19	0.42
1:B:20:ILE:HG21	1:B:372:ILE:HD11	2.01	0.42
1:B:417:ALA:HB1	1:B:484:ARG:O	2.19	0.42
1:C:371:GLN:CB	1:C:396:VAL:HG22	2.48	0.42
1:C:408:ALA:CB	1:C:463:THR:HG22	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:LEU:HD22	1:A:13:LEU:HD21	2.01	0.42
1:A:29:LEU:HA	1:A:29:LEU:HD23	1.83	0.42
1:A:503:GLN:HG3	1:A:504:GLY:O	2.19	0.42
1:C:360:ILE:HD13	1:C:360:ILE:HA	1.90	0.42
1:D:288:ARG:HA	1:D:291:LEU:HB3	2.02	0.42
1:A:11:LEU:O	1:A:13:LEU:HG	2.20	0.42
1:A:24:ASN:ND2	1:A:381:GLY:O	2.49	0.42
1:B:283:THR:CG2	1:B:284:LEU:N	2.83	0.42
1:B:414:PHE:O	1:B:416:THR:N	2.52	0.42
1:B:422:PHE:CD2	1:B:501:TRP:CH2	3.07	0.42
1:B:484:ARG:HD2	1:B:497:GLU:O	2.19	0.42
1:D:318:PHE:CD1	1:D:367:HIS:ND1	2.86	0.42
1:A:379:GLN:HG2	1:B:438:TRP:NE1	2.34	0.42
1:B:108:PRO:HG2	1:B:264:ASN:HD21	1.85	0.42
1:B:143:LEU:HD11	1:B:208:PHE:HD2	1.85	0.42
1:C:20:ILE:O	1:C:393:GLU:HA	2.20	0.42
1:C:66:TRP:CD1	1:C:66:TRP:H	2.38	0.42
1:D:266:ARG:NE	1:D:266:ARG:CA	2.76	0.42
1:D:212:THR:HA	1:D:213:PRO:HD2	1.73	0.42
1:A:407:CYS:HB3	1:A:464:VAL:HG12	2.02	0.42
1:D:436:ILE:HG12	1:D:457:ARG:NH1	2.34	0.42
1:B:125:LEU:HA	1:B:125:LEU:HD13	1.65	0.41
1:D:421:TRP:CH2	1:D:481:CYS:HB3	2.54	0.41
1:D:210:ASN:OD1	1:D:210:ASN:N	2.53	0.41
1:D:318:PHE:HB3	1:D:367:HIS:ND1	2.35	0.41
1:D:321:THR:HG22	1:D:322:ASP:N	2.36	0.41
1:C:370:TYR:CD2	1:C:370:TYR:C	2.98	0.41
1:D:293:LEU:O	1:D:297:MET:HG3	2.20	0.41
1:A:474:GLU:CD	1:A:474:GLU:C	2.88	0.41
1:B:7:SER:CB	1:C:101:GLN:OE1	2.68	0.41
1:B:367:HIS:CE1	1:B:396:VAL:HG13	2.55	0.41
1:D:10:GLY:C	1:D:403:LEU:HD23	2.44	0.41
1:D:440:VAL:HG13	1:D:442:SER:H	1.85	0.41
1:D:441:GLY:HA2	1:D:457:ARG:HH12	1.86	0.41
1:A:65:SER:O	1:A:68:THR:HG22	2.20	0.41
1:C:45:PRO:HB2	1:C:301:VAL:HG13	2.02	0.41
1:A:150:GLU:HA	1:A:153:PHE:HB3	2.03	0.41
1:C:369:ALA:O	1:C:370:TYR:C	2.63	0.41
1:D:436:ILE:O	1:D:436:ILE:HG23	2.20	0.41
1:B:215:ILE:O	1:B:219:GLN:N	2.54	0.41
1:D:86:ALA:HA	1:D:89:MET:HE2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:LEU:CD2	1:B:401:SER:HB2	2.51	0.41
1:B:97:VAL:HA	1:B:100:THR:HG22	2.03	0.41
1:B:283:THR:HG23	1:B:382:LEU:HD21	2.01	0.41
1:A:217:GLN:O	1:A:221:LEU:HD22	2.21	0.41
1:A:287:LEU:HD12	1:A:382:LEU:HD22	2.03	0.41
1:B:353:ASP:OD1	1:B:354:LYS:N	2.51	0.40
1:C:82:GLU:OE1	1:C:82:GLU:N	2.53	0.40
1:C:192:ALA:O	1:C:288:ARG:HD3	2.21	0.40
1:B:108:PRO:HG2	1:B:264:ASN:ND2	2.37	0.40
1:B:143:LEU:HD23	1:B:143:LEU:HA	1.67	0.40
1:A:492:PRO:HB2	1:A:498:TYR:HE1	1.86	0.40
1:D:262:ASN:O	1:D:266:ARG:HG2	2.21	0.40
1:A:63:THR:N	1:A:64:PRO:HD2	2.36	0.40
1:C:274:THR:HA	1:C:280:TYR:CG	2.57	0.40
1:C:421:TRP:CH2	1:C:481:CYS:HB3	2.56	0.40

There are no symmetry-related clashes.

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	398/521 (76%)	378 (95%)	20 (5%)	0	100	100
1	B	415/521 (80%)	390 (94%)	24 (6%)	1 (0%)	43	71
1	C	394/521 (76%)	374 (95%)	20 (5%)	0	100	100
1	D	369/521 (71%)	350 (95%)	17 (5%)	2 (0%)	24	55
All	All	1576/2084 (76%)	1492 (95%)	81 (5%)	3 (0%)	43	71

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	213	PRO
1	B	10	GLY
1	D	356	VAL

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	335/427 (78%)	332 (99%)	3 (1%)	70	77
1	B	347/427 (81%)	342 (99%)	5 (1%)	59	73
1	C	331/427 (78%)	331 (100%)	0	100	100
1	D	310/427 (73%)	308 (99%)	2 (1%)	78	81
All	All	1323/1708 (78%)	1313 (99%)	10 (1%)	73	78

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	414	PHE
1	A	457	ARG
1	A	475	ASP
1	B	6	GLU
1	B	16	THR
1	B	17	THR
1	B	293	LEU
1	B	321	THR
1	D	416	THR
1	D	437	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	ASN
1	A	189	ASN
1	B	148	GLN
1	B	313	ASN

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Mol	Chain	Res	Type
1	B	315	GLN
1	C	27	GLN
1	C	28	ASN
1	C	91	ASN
1	C	148	GLN
1	D	262	ASN
1	D	371	GLN
1	D	379	GLN
1	D	462	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 ⓘ

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	412/521 (79%)	0.11	14 (3%)	48	28	66, 123, 176, 241	0
1	B	429/521 (82%)	0.03	16 (3%)	45	26	57, 103, 144, 163	0
1	C	412/521 (79%)	0.10	16 (3%)	43	25	49, 99, 153, 182	0
1	D	387/521 (74%)	0.24	26 (6%)	24	13	69, 129, 181, 206	0
All	All	1640/2084 (78%)	0.12	72 (4%)	39	22	49, 112, 168, 241	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	486	ARG	6.3
1	D	489	THR	5.3
1	D	215	ILE	3.9
1	A	365	LYS	3.9
1	C	287	LEU	3.9
1	D	216	ASN	3.7
1	C	364	GLU	3.6
1	B	111	ILE	3.6
1	D	212	THR	3.6
1	B	112	THR	3.5
1	B	489	THR	3.4
1	A	376	ALA	3.1
1	B	16	THR	3.1
1	B	207	ASP	3.1
1	B	490	LEU	3.1
1	D	161	LEU	3.1
1	D	467	GLN	3.1
1	D	32	GLN	3.1
1	A	467	GLN	3.0
1	B	271	ALA	2.9
1	C	13	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	64	PRO	2.9
1	D	402	SER	2.8
1	C	125	LEU	2.8
1	C	19	VAL	2.7
1	B	10	GLY	2.7
1	D	14	LYS	2.7
1	D	97	VAL	2.7
1	A	29	LEU	2.7
1	B	278	PRO	2.6
1	D	469	ASP	2.6
1	D	9	GLY	2.6
1	A	22	THR	2.6
1	B	125	LEU	2.6
1	A	140	ALA	2.6
1	C	9	GLY	2.5
1	A	499	ASP	2.5
1	B	297	MET	2.4
1	C	111	ILE	2.3
1	D	69	ALA	2.3
1	D	21	ASP	2.3
1	B	15	THR	2.3
1	D	405	LEU	2.3
1	C	442	SER	2.3
1	B	299	TYR	2.3
1	C	210	ASN	2.3
1	D	36	ILE	2.2
1	D	49	ALA	2.2
1	A	32	GLN	2.2
1	B	443	THR	2.2
1	D	314	ASN	2.2
1	A	373	LEU	2.2
1	A	378	LYS	2.2
1	D	12	VAL	2.1
1	C	295	ASN	2.1
1	D	213	PRO	2.1
1	A	222	ALA	2.1
1	C	289	SER	2.1
1	D	502	GLY	2.1
1	C	314	ASN	2.1
1	A	25	ASP	2.1
1	C	90	ILE	2.1
1	C	128	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	457	ARG	2.1
1	D	488	PHE	2.1
1	D	332	CYS	2.1
1	C	286	ALA	2.1
1	D	369	ALA	2.1
1	B	17	THR	2.0
1	A	368	GLU	2.0
1	B	132	MET	2.0
1	D	305	GLY	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.