



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:09 AM UTC

PDB ID : 1KCU / pdb_00001kcu
Title : CRYSTAL STRUCTURE OF ANTIBODY PC287
Authors : Nair, D.T.; Singh, K.; Siddiqui, Z.; Nayak, B.P.; Rao, K.V.S.; Salunke, D.M.
Deposited on : 2001-11-11
Resolution : 2.20 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

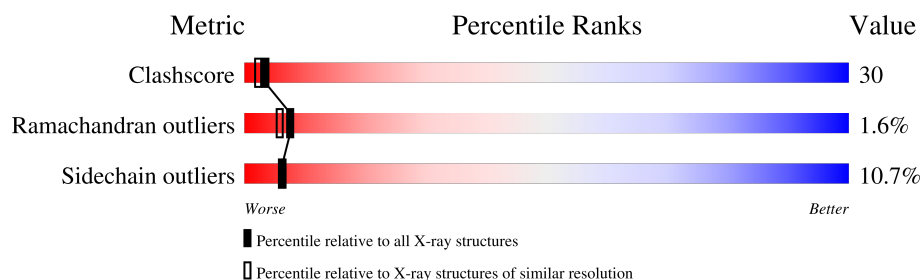
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	214	
2	H	217	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3392 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 PC287 IMMUNOGLOBULIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	214	Total	C	N	O	S	0	0	0
			1648	1023	275	342	8			

- Molecule 2 is a protein called PC287 IMMUNOGLOBULIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	217	Total	C	N	O	S	0	0	0
			1635	1040	266	323	6			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	59	Total	O	0	0
			59	59		
3	H	50	Total	O	0	0
			50	50		

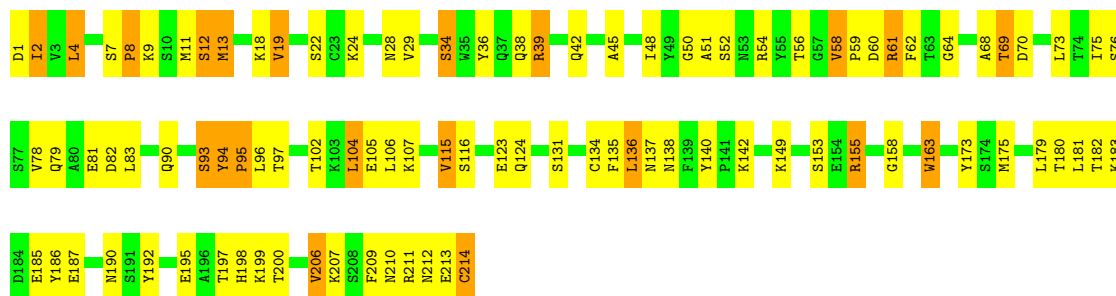
3 Residue-property plots [i](#)

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. 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. 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.

Note EDS was not executed.

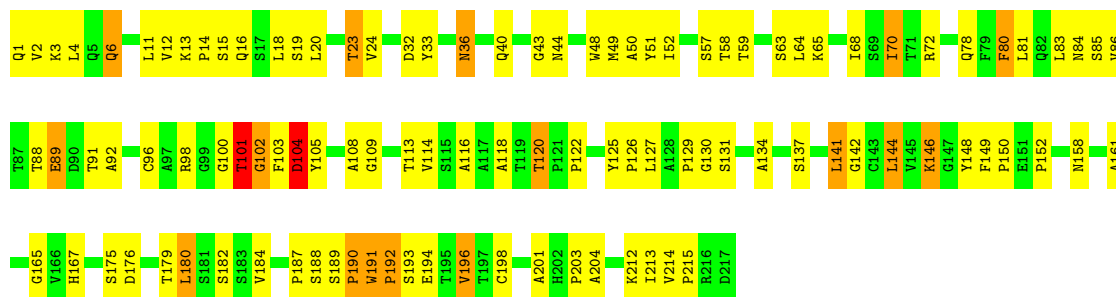
• Molecule 1: PC287 IMMUNOGLOBULIN

Chain L: 



• Molecule 2: PC287 IMMUNOGLOBULIN

Chain H: 



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	73.76Å 71.64Å 87.72Å 90.00° 98.14° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20	Depositor
% Data completeness (in resolution range)	85.0 (50.00-2.20)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.5	Depositor
R, R_{free}	0.200 , 0.249	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3392	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

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	L	0.62	2/1685 (0.1%)	1.14	15/2290 (0.7%)
2	H	0.64	1/1681 (0.1%)	1.55	19/2307 (0.8%)
All	All	0.63	3/3366 (0.1%)	1.36	34/4597 (0.7%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	94	TYR	CA-CB	-6.20	1.43	1.53
1	L	94	TYR	CB-CG	-6.00	1.38	1.51
2	H	89	GLU	CA-C	5.38	1.60	1.52

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	191	TRP	CA-C-N	32.44	160.40	119.84
2	H	191	TRP	C-N-CA	32.44	160.40	119.84
2	H	191	TRP	C-N-CD	-14.22	66.72	125.00
2	H	191	TRP	CA-C-O	10.94	126.40	119.29
2	H	191	TRP	N-CA-C	10.16	126.66	108.47
1	L	79	GLN	OE1-CD-NE2	-9.91	112.69	122.60
1	L	94	TYR	CA-C-N	-9.39	108.10	119.84
1	L	94	TYR	C-N-CA	-9.39	108.10	119.84
2	H	104	ASP	CA-CB-CG	9.35	121.95	112.60
2	H	101	THR	N-CA-C	-8.04	95.77	108.63
2	H	191	TRP	CA-CB-CG	-7.53	99.30	113.60
1	L	79	GLN	CG-CD-NE2	6.74	126.51	116.40
2	H	191	TRP	O-C-N	6.65	126.61	121.23
2	H	100	GLY	O-C-N	6.49	131.14	122.70
1	L	58	VAL	N-CA-C	6.46	114.19	108.63
1	L	137	ASN	N-CA-C	6.37	120.30	110.42
1	L	94	TYR	CB-CG-CD2	-6.24	111.44	120.80
2	H	36	ASN	N-CA-C	6.16	119.68	110.14
1	L	182	THR	N-CA-C	-6.07	102.35	110.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	19	VAL	CB-CA-C	-6.02	103.51	111.70
2	H	104	ASP	CA-C-N	-5.96	113.77	122.65
2	H	104	ASP	C-N-CA	-5.96	113.77	122.65
1	L	206	VAL	N-CA-C	5.94	116.68	108.12
1	L	93	SER	CA-C-N	-5.81	107.62	121.80
1	L	93	SER	C-N-CA	-5.81	107.62	121.80
1	L	94	TYR	N-CA-C	5.81	122.65	109.81
2	H	108	ALA	N-CA-C	-5.50	106.52	113.18
2	H	146	LYS	N-CA-C	5.47	118.62	110.14
1	L	138	ASN	N-CA-C	5.32	118.67	111.17
2	H	191	TRP	CB-CA-C	-5.27	104.02	110.08
2	H	96	CYS	N-CA-C	-5.17	101.45	109.72
2	H	120	THR	CA-C-N	5.07	123.36	119.66
2	H	120	THR	C-N-CA	5.07	123.36	119.66
1	L	115	VAL	N-CA-C	5.04	116.07	109.21

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	L	1648	0	1569	96	0
2	H	1635	0	1599	101	0
3	H	50	0	0	6	0
3	L	59	0	0	5	0
All	All	3392	0	3168	193	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 30.

All (193) 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:L:68:ALA:O	1:L:69:THR:HB	1.52	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:189:SER:HB2	2:H:190:PRO:CD	1.93	0.97
2:H:101:THR:O	2:H:102:GLY:O	1.82	0.97
1:L:7:SER:HB3	1:L:8:PRO:HD2	1.50	0.94
2:H:12:VAL:HG11	2:H:18:LEU:HD23	1.54	0.90
1:L:28:ASN:ND2	1:L:68:ALA:HB1	1.87	0.89
2:H:91:THR:HG22	2:H:114:VAL:H	1.37	0.89
1:L:68:ALA:O	1:L:69:THR:CB	2.18	0.88
1:L:38:GLN:HE22	2:H:40:GLN:HE22	1.22	0.88
2:H:118:ALA:O	2:H:120:THR:HG23	1.74	0.87
1:L:39:ARG:HH11	1:L:39:ARG:HB2	1.42	0.84
2:H:189:SER:HB2	2:H:190:PRO:HD3	1.59	0.81
1:L:7:SER:HB3	1:L:8:PRO:CD	2.11	0.80
2:H:23:THR:HG22	3:H:266:HOH:O	1.80	0.80
1:L:7:SER:CB	1:L:8:PRO:CD	2.60	0.79
2:H:52:ILE:HB	2:H:70:ILE:HG21	1.67	0.77
2:H:189:SER:HB2	2:H:190:PRO:HD2	1.68	0.76
1:L:4:LEU:HD11	1:L:90:GLN:CG	2.17	0.75
2:H:23:THR:CG2	3:H:266:HOH:O	2.35	0.74
2:H:68:ILE:HG22	2:H:83:LEU:HD23	1.68	0.74
2:H:88:THR:CG2	2:H:89:GLU:OE2	2.36	0.74
1:L:94:TYR:O	1:L:96:LEU:N	2.22	0.72
1:L:13:MET:HE2	1:L:104:LEU:HG	1.72	0.72
2:H:6:GLN:H	2:H:6:GLN:NE2	1.87	0.72
2:H:187:PRO:HG2	2:H:190:PRO:HD2	1.70	0.71
1:L:95:PRO:O	1:L:97:THR:HG23	1.90	0.71
2:H:23:THR:HB	2:H:78:GLN:HE21	1.56	0.71
2:H:189:SER:CB	2:H:190:PRO:CD	2.68	0.71
1:L:2:ILE:HD11	1:L:90:GLN:HG2	1.73	0.70
1:L:48:ILE:HD13	1:L:54:ARG:HA	1.73	0.70
2:H:40:GLN:O	2:H:92:ALA:HB1	1.91	0.70
1:L:2:ILE:HG13	1:L:2:ILE:O	1.91	0.69
2:H:88:THR:O	2:H:91:THR:HG23	1.93	0.69
2:H:101:THR:O	2:H:101:THR:OG1	2.07	0.68
2:H:49:MET:HE1	2:H:81:LEU:CD2	2.24	0.68
2:H:214:VAL:HG13	2:H:215:PRO:HD2	1.75	0.68
1:L:38:GLN:HA	3:L:225:HOH:O	1.93	0.67
1:L:61:ARG:HA	1:L:76:SER:OG	1.93	0.67
2:H:36:ASN:OD1	2:H:51:TYR:HB3	1.94	0.67
1:L:83:LEU:HD21	1:L:106:LEU:HB2	1.75	0.67
2:H:129:PRO:HD3	2:H:141:LEU:HD12	1.78	0.66
1:L:210:ASN:ND2	3:L:222:HOH:O	2.29	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:191:TRP:CD1	2:H:196:VAL:HG13	2.32	0.64
2:H:88:THR:HG23	2:H:89:GLU:OE2	1.97	0.64
1:L:61:ARG:O	1:L:75:ILE:HA	1.98	0.64
2:H:51:TYR:C	2:H:51:TYR:CD1	2.75	0.64
2:H:122:PRO:HB3	2:H:148:TYR:HB3	1.80	0.63
1:L:95:PRO:O	1:L:96:LEU:C	2.40	0.62
2:H:191:TRP:HD1	2:H:196:VAL:HG13	1.65	0.62
1:L:158:GLY:O	1:L:179:LEU:HA	1.99	0.62
2:H:16:GLN:O	2:H:86:VAL:HG22	2.01	0.60
2:H:6:GLN:H	2:H:6:GLN:HE21	1.49	0.58
1:L:4:LEU:HD11	1:L:90:GLN:HG3	1.83	0.58
1:L:7:SER:OG	1:L:8:PRO:HD3	2.04	0.58
2:H:101:THR:O	2:H:102:GLY:C	2.47	0.57
2:H:14:PRO:O	2:H:15:SER:HB2	2.03	0.57
1:L:173:TYR:HB3	3:L:261:HOH:O	2.04	0.57
2:H:11:LEU:HD22	2:H:113:THR:O	2.05	0.57
2:H:49:MET:CE	2:H:81:LEU:CD2	2.83	0.57
1:L:2:ILE:HD13	1:L:29:VAL:HG11	1.86	0.57
2:H:19:SER:HA	2:H:81:LEU:O	2.05	0.56
2:H:187:PRO:HG2	2:H:190:PRO:HG2	1.88	0.56
1:L:48:ILE:CD1	1:L:54:ARG:HA	2.36	0.56
2:H:52:ILE:HB	2:H:70:ILE:CG2	2.35	0.56
2:H:1:GLN:HG3	2:H:2:VAL:H	1.72	0.55
1:L:142:LYS:HG2	1:L:173:TYR:CZ	2.42	0.55
2:H:180:LEU:C	2:H:180:LEU:HD23	2.32	0.54
2:H:175:SER:OG	2:H:176:ASP:N	2.40	0.54
2:H:11:LEU:HD21	2:H:149:PHE:HZ	1.72	0.54
2:H:104:ASP:HB2	2:H:105:TYR:CD2	2.43	0.54
1:L:28:ASN:HD21	1:L:68:ALA:HB1	1.70	0.54
2:H:101:THR:C	2:H:102:GLY:O	2.48	0.53
1:L:190:ASN:HD21	1:L:212:ASN:H	1.55	0.53
2:H:127:LEU:HB2	2:H:142:GLY:C	2.34	0.53
2:H:91:THR:CG2	2:H:114:VAL:H	2.15	0.53
1:L:124:GLN:HE22	1:L:131:SER:H	1.56	0.53
2:H:49:MET:CE	2:H:81:LEU:HD22	2.39	0.53
1:L:142:LYS:HG2	1:L:173:TYR:CE1	2.44	0.53
2:H:11:LEU:HD23	2:H:113:THR:HB	1.90	0.53
1:L:2:ILE:CD1	1:L:90:GLN:HG2	2.38	0.52
1:L:68:ALA:C	1:L:69:THR:HG22	2.34	0.52
2:H:187:PRO:HG2	2:H:190:PRO:CD	2.38	0.52
2:H:149:PHE:CE2	2:H:150:PRO:HB3	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:60:ASP:C	1:L:62:PHE:H	2.17	0.52
1:L:2:ILE:HD13	1:L:29:VAL:CG1	2.39	0.52
1:L:54:ARG:HD3	1:L:59:PRO:O	2.10	0.52
1:L:81:GLU:H	1:L:81:GLU:CD	2.18	0.51
2:H:63:SER:C	2:H:64:LEU:HD12	2.35	0.51
1:L:163:TRP:N	1:L:163:TRP:CD1	2.78	0.51
1:L:24:LYS:HB3	1:L:24:LYS:NZ	2.26	0.51
1:L:8:PRO:HB3	1:L:11:MET:HE3	1.92	0.51
1:L:4:LEU:CD1	1:L:90:GLN:HG3	2.39	0.51
1:L:206:VAL:HB	3:L:262:HOH:O	2.10	0.51
2:H:88:THR:HG22	2:H:89:GLU:OE2	2.09	0.51
1:L:39:ARG:HH11	1:L:39:ARG:CB	2.19	0.51
1:L:13:MET:HE3	1:L:105:GLU:O	2.11	0.50
2:H:146:LYS:HB3	2:H:179:THR:HG23	1.92	0.50
1:L:7:SER:HB3	1:L:22:SER:HB3	1.93	0.50
2:H:129:PRO:HD3	2:H:141:LEU:CD1	2.42	0.50
1:L:149:LYS:NZ	1:L:195:GLU:CD	2.70	0.50
2:H:52:ILE:HD13	2:H:72:ARG:HD3	1.92	0.50
2:H:130:GLY:O	2:H:131:SER:HB3	2.12	0.50
1:L:2:ILE:HD11	1:L:4:LEU:HD11	1.94	0.49
2:H:189:SER:CB	2:H:190:PRO:HD3	2.35	0.49
1:L:149:LYS:NZ	1:L:195:GLU:OE1	2.45	0.49
2:H:6:GLN:NE2	2:H:109:GLY:H	2.10	0.49
1:L:195:GLU:HG3	1:L:206:VAL:HG12	1.94	0.49
2:H:158:ASN:HB3	2:H:161:ALA:HB3	1.95	0.49
1:L:7:SER:CB	1:L:8:PRO:HD3	2.40	0.49
2:H:165:GLY:O	2:H:184:VAL:HA	2.13	0.49
2:H:49:MET:HE1	2:H:81:LEU:HD21	1.91	0.48
2:H:13:LYS:HD2	2:H:116:ALA:HA	1.95	0.48
1:L:180:THR:O	1:L:181:LEU:HD23	2.13	0.48
2:H:80:PHE:CD1	2:H:80:PHE:N	2.80	0.48
1:L:61:ARG:NH1	1:L:82:ASP:OD2	2.47	0.48
1:L:94:TYR:HE1	2:H:51:TYR:HE2	1.62	0.48
1:L:155:ARG:HH11	1:L:155:ARG:HG3	1.79	0.48
1:L:116:SER:O	1:L:134:CYS:HA	2.14	0.48
2:H:51:TYR:C	2:H:51:TYR:HD1	2.21	0.47
1:L:8:PRO:O	1:L:102:THR:HG23	2.14	0.47
1:L:124:GLN:HG3	2:H:125:TYR:CZ	2.49	0.47
1:L:54:ARG:HD3	1:L:58:VAL:HG12	1.96	0.47
2:H:84:ASN:O	2:H:85:SER:C	2.58	0.47
2:H:201:ALA:O	2:H:203:PRO:HD3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:198:HIS:CD2	1:L:200:THR:HG23	2.49	0.47
1:L:90:GLN:CD	1:L:90:GLN:C	2.84	0.46
2:H:48:TRP:HH2	2:H:59:THR:HG22	1.81	0.46
2:H:129:PRO:C	2:H:131:SER:H	2.22	0.46
1:L:13:MET:SD	1:L:78:VAL:HG11	2.55	0.46
2:H:187:PRO:C	2:H:189:SER:H	2.23	0.46
1:L:28:ASN:ND2	1:L:68:ALA:CB	2.69	0.46
1:L:1:ASP:CG	1:L:2:ILE:H	2.23	0.46
1:L:24:LYS:HB3	1:L:24:LYS:HZ3	1.80	0.46
1:L:185:GLU:OE2	1:L:185:GLU:HA	2.16	0.45
2:H:91:THR:HG22	2:H:114:VAL:N	2.18	0.45
1:L:213:GLU:O	1:L:214:CYS:C	2.59	0.45
2:H:84:ASN:HB3	3:H:255:HOH:O	2.16	0.45
1:L:60:ASP:O	1:L:62:PHE:N	2.49	0.45
2:H:1:GLN:HG3	2:H:2:VAL:N	2.30	0.45
2:H:64:LEU:HD12	2:H:64:LEU:N	2.32	0.45
1:L:60:ASP:C	1:L:62:PHE:N	2.74	0.45
1:L:186:TYR:CE2	1:L:211:ARG:HD3	2.52	0.45
2:H:84:ASN:CB	3:H:255:HOH:O	2.65	0.45
2:H:48:TRP:CZ2	2:H:50:ALA:HA	2.52	0.45
2:H:191:TRP:HZ2	2:H:213:ILE:O	1.99	0.44
1:L:2:ILE:CG1	1:L:90:GLN:HG2	2.47	0.44
2:H:52:ILE:HG13	2:H:58:THR:HG22	2.00	0.44
2:H:49:MET:CE	2:H:81:LEU:HD21	2.48	0.44
2:H:187:PRO:HG2	2:H:190:PRO:CG	2.48	0.44
1:L:136:LEU:N	1:L:136:LEU:HD22	2.33	0.44
2:H:213:ILE:HA	3:H:229:HOH:O	2.17	0.44
2:H:167:HIS:O	2:H:182:SER:HA	2.18	0.44
1:L:207:LYS:HD3	1:L:207:LYS:HA	1.88	0.44
1:L:4:LEU:HD11	1:L:90:GLN:HG2	1.98	0.43
1:L:136:LEU:HB2	1:L:175:MET:HG2	2.01	0.43
2:H:125:TYR:HD2	2:H:144:LEU:HD12	1.82	0.43
1:L:28:ASN:HD21	1:L:68:ALA:CB	2.29	0.43
1:L:107:LYS:HA	1:L:140:TYR:OH	2.19	0.43
2:H:192:PRO:HG2	2:H:193:SER:H	1.83	0.43
2:H:192:PRO:CG	2:H:193:SER:H	2.31	0.43
1:L:192:TYR:HB2	1:L:209:PHE:CE2	2.54	0.42
2:H:49:MET:HE3	2:H:81:LEU:HD22	2.00	0.42
2:H:126:PRO:HB2	3:H:229:HOH:O	2.18	0.42
1:L:183:LYS:HE2	1:L:187:GLU:OE2	2.18	0.42
2:H:63:SER:OG	2:H:64:LEU:HD12	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:158:ASN:OD1	2:H:196:VAL:HA	2.19	0.42
2:H:23:THR:HB	2:H:78:GLN:NE2	2.28	0.42
2:H:32:ASP:CG	2:H:33:TYR:H	2.28	0.42
1:L:96:LEU:HD22	2:H:103:PHE:HZ	1.85	0.42
1:L:34:SER:HA	1:L:48:ILE:O	2.20	0.42
1:L:1:ASP:HB2	1:L:95:PRO:HB2	2.01	0.42
1:L:149:LYS:HA	1:L:153:SER:O	2.20	0.42
2:H:43:GLY:O	2:H:44:ASN:HB2	2.20	0.42
1:L:135:PHE:C	1:L:136:LEU:HD22	2.46	0.41
1:L:210:ASN:ND2	3:L:269:HOH:O	2.53	0.41
1:L:115:VAL:HG22	1:L:136:LEU:HD13	2.02	0.41
1:L:52:SER:HA	1:L:64:GLY:O	2.19	0.41
1:L:211:ARG:C	1:L:213:GLU:H	2.29	0.41
1:L:36:TYR:HA	1:L:45:ALA:O	2.21	0.41
1:L:69:THR:HG23	1:L:70:ASP:OD1	2.21	0.41
1:L:190:ASN:HD21	1:L:212:ASN:N	2.19	0.41
2:H:11:LEU:HD13	2:H:11:LEU:C	2.45	0.41
2:H:150:PRO:HD2	2:H:204:ALA:CB	2.51	0.41
2:H:3:LYS:C	2:H:4:LEU:HD12	2.46	0.41
1:L:12:SER:HG	1:L:105:GLU:CD	2.29	0.40
1:L:50:GLY:O	1:L:51:ALA:HB3	2.21	0.40
1:L:190:ASN:O	1:L:210:ASN:HA	2.22	0.40
2:H:212:LYS:HB3	2:H:212:LYS:HE2	1.83	0.40
1:L:2:ILE:CD1	1:L:29:VAL:HG11	2.50	0.40
1:L:1:ASP:CG	1:L:2:ILE:N	2.80	0.40
2:H:104:ASP:HB2	2:H:105:TYR:HD2	1.83	0.40
2:H:91:THR:HG22	2:H:114:VAL:HB	2.02	0.40
2:H:137:SER:O	2:H:188:SER:N	2.50	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	L	212/214 (99%)	193 (91%)	16 (8%)	3 (1%)	9	7
2	H	215/217 (99%)	184 (86%)	27 (13%)	4 (2%)	6	4
All	All	427/431 (99%)	377 (88%)	43 (10%)	7 (2%)	7	6

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	69	THR
2	H	102	GLY
2	H	192	PRO
1	L	61	ARG
2	H	190	PRO
2	H	134	ALA
1	L	95	PRO

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	L	188/188 (100%)	166 (88%)	22 (12%)	5	5
2	H	185/185 (100%)	167 (90%)	18 (10%)	8	8
All	All	373/373 (100%)	333 (89%)	40 (11%)	6	6

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	2	ILE
1	L	4	LEU
1	L	8	PRO
1	L	9	LYS
1	L	12	SER
1	L	13	MET
1	L	18	LYS
1	L	19	VAL
1	L	34	SER

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Mol	Chain	Res	Type
1	L	39	ARG
1	L	42	GLN
1	L	56	THR
1	L	73	LEU
1	L	93	SER
1	L	104	LEU
1	L	123	GLU
1	L	136	LEU
1	L	155	ARG
1	L	163	TRP
1	L	197	THR
1	L	199	LYS
1	L	214	CYS
2	H	6	GLN
2	H	20	LEU
2	H	23	THR
2	H	24	VAL
2	H	57	SER
2	H	65	LYS
2	H	70	ILE
2	H	80	PHE
2	H	98	ARG
2	H	101	THR
2	H	104	ASP
2	H	141	LEU
2	H	144	LEU
2	H	152	PRO
2	H	180	LEU
2	H	194	GLU
2	H	196	VAL
2	H	198	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	28	ASN
1	L	37	GLN
1	L	42	GLN
1	L	87	HIS
1	L	124	GLN
1	L	198	HIS
1	L	210	ASN

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Mol	Chain	Res	Type
2	H	6	GLN
2	H	40	GLN
2	H	78	GLN
2	H	167	HIS
2	H	174	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](#)

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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.