



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:12 PM UTC

PDB ID : 3HR5 / pdb_00003hr5
Title : M1prime peptide from IgE bound by humanized antibody 47H4 Fab
Authors : Eigenbrot, C.W.; Ultsch, M.H.
Deposited on : 2009-06-08
Resolution : 2.40 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

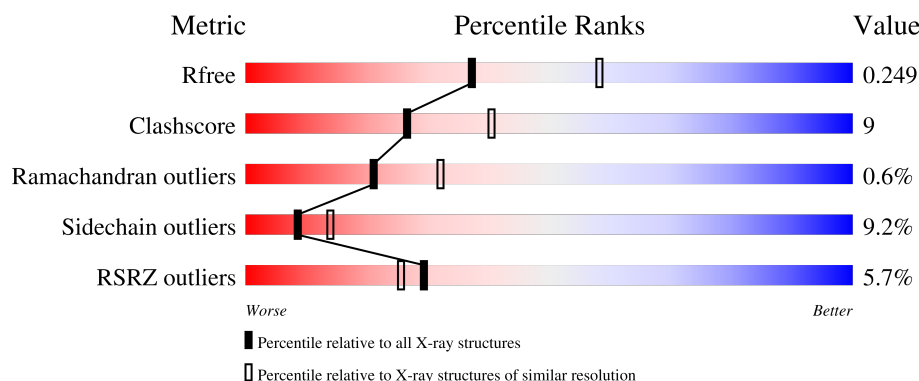
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	219	4% 71% 24% .
1	L	219	3% 75% 21% .
1	P	219	5% 77% 20% .
1	Q	219	5% 78% 19% .
2	B	226	6% 79% 17% ..

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Mol	Chain	Length	Quality of chain
2	H	226	
2	I	226	
2	J	226	
3	R	35	
3	S	35	
3	T	35	
3	V	35	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GOL	P	402	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 14144 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 Fab h47H4 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	219	Total	C	N	O	S	0	1	0
			1702	1063	290	342	7			
1	A	219	Total	C	N	O	S	0	2	0
			1713	1069	294	343	7			
1	P	219	Total	C	N	O	S	0	3	0
			1722	1074	295	346	7			
1	Q	219	Total	C	N	O	S	0	4	0
			1726	1077	295	347	7			

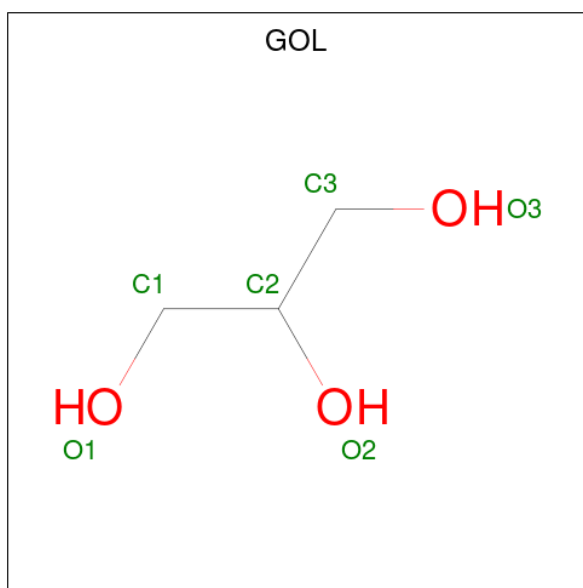
- Molecule 2 is a protein called Fab h47H4 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	220	Total	C	N	O	S	0	2	0
			1662	1049	274	331	8			
2	B	220	Total	C	N	O	S	0	1	0
			1656	1046	273	330	7			
2	I	220	Total	C	N	O	S	0	0	0
			1648	1042	272	327	7			
2	J	220	Total	C	N	O	S	0	2	0
			1660	1048	274	330	8			

- Molecule 3 is a protein called M1prime-derived peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	R	9	Total	C	N	O	0	0	0
			71	40	17	14			
3	S	11	Total	C	N	O	0	0	0
			84	48	19	17			
3	T	11	Total	C	N	O	0	0	0
			86	51	19	16			
3	V	10	Total	C	N	O	0	0	0
			78	45	18	15			

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	P	1	Total	C	O	0	0
			6	3	3		
4	Q	1	Total	C	O	0	0
			6	3	3		
4	J	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	30	Total	O	0	0
			30	30		
5	H	49	Total	O	0	0
			49	49		
5	A	45	Total	O	0	0
			45	45		
5	B	53	Total	O	0	0
			53	53		
5	P	49	Total	O	0	0
			49	49		
5	I	15	Total	O	0	0
			15	15		
5	Q	35	Total	O	0	0
			35	35		

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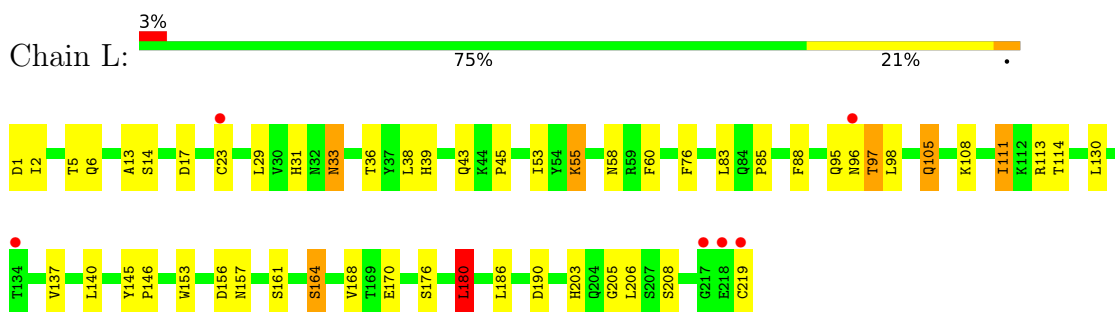
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	J	35	Total 35	O 35	0	0
5	S	1	Total 1	O 1	0	0

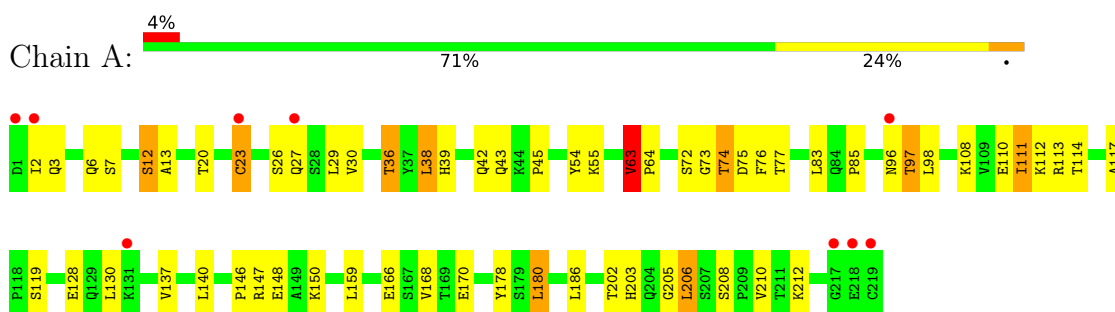
3 Residue-property plots [i](#)

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.

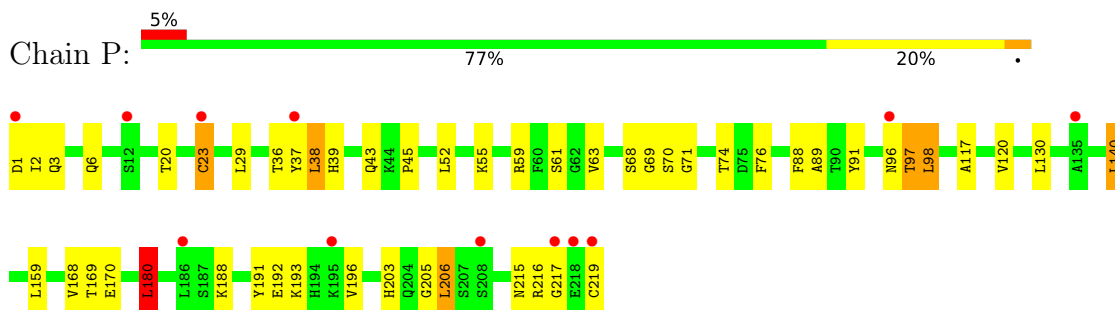
- Molecule 1: Fab h47H4 light chain



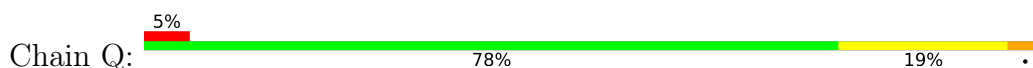
- Molecule 1: Fab h47H4 light chain

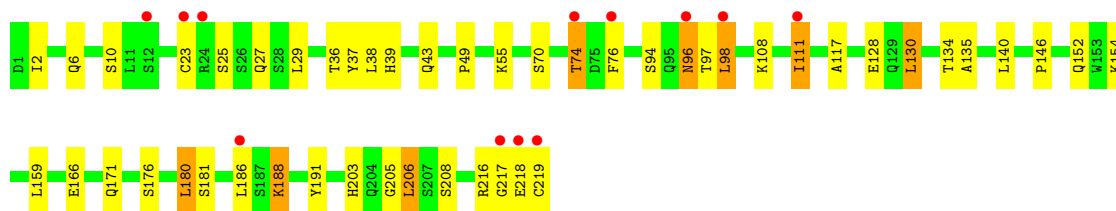


- Molecule 1: Fab h47H4 light chain

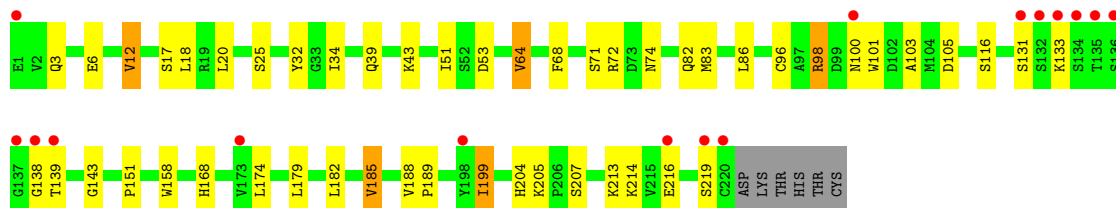
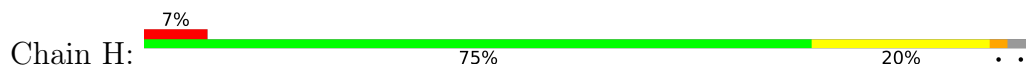


- Molecule 1: Fab h47H4 light chain

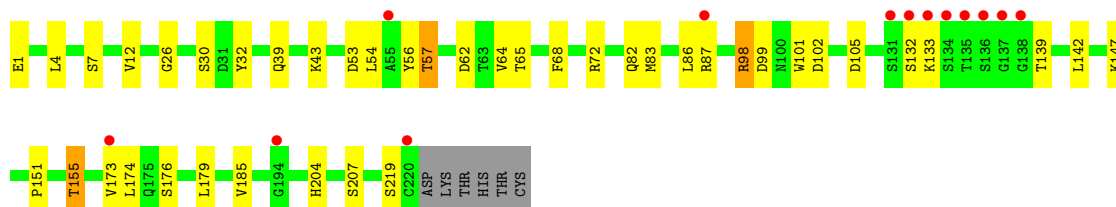
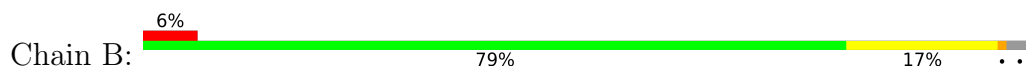




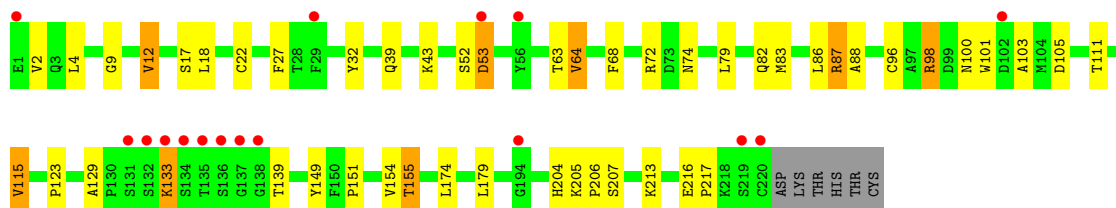
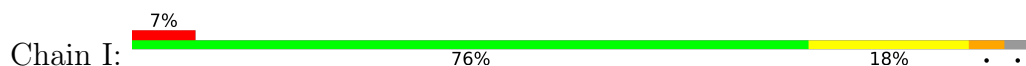
• Molecule 2: Fab h47H4 heavy chain



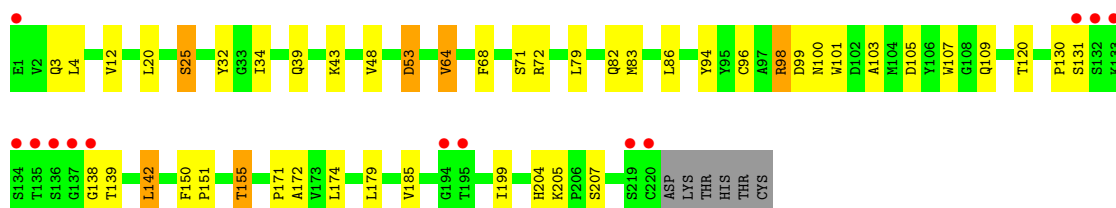
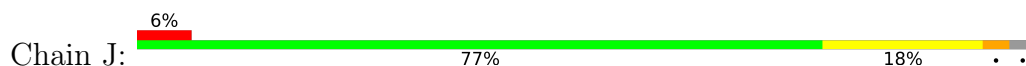
• Molecule 2: Fab h47H4 heavy chain



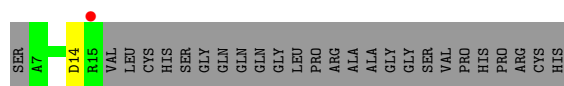
• Molecule 2: Fab h47H4 heavy chain



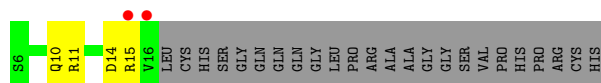
• Molecule 2: Fab h47H4 heavy chain



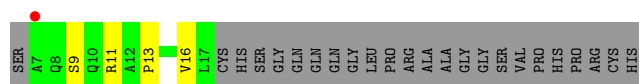
- Molecule 3: M1prime-derived peptide



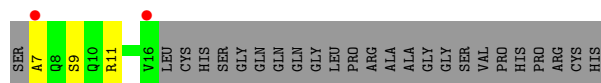
- Molecule 3: M1prime-derived peptide



- Molecule 3: M1prime-derived peptide



- Molecule 3: M1prime-derived peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.76Å 183.82Å 194.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 50.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	89.0 (50.00-2.40) 88.9 (50.00-2.40)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.213 , 0.253 0.210 , 0.249	Depositor DCC
R_{free} test set	1040 reflections (1.27%)	wwPDB-VP
Wilson B-factor (Å ²)	44.2	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.5	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	0.93	EDS
Total number of atoms	14144	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.87% 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:
GOL

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	0.63	0/1751	0.86	1/2376 (0.0%)
1	L	0.58	0/1740	0.82	1/2362 (0.0%)
1	P	0.59	0/1760	0.82	1/2388 (0.0%)
1	Q	0.59	0/1764	0.85	1/2393 (0.0%)
2	B	0.65	0/1696	0.89	2/2315 (0.1%)
2	H	0.65	0/1702	0.86	1/2323 (0.0%)
2	I	0.60	0/1688	0.85	2/2304 (0.1%)
2	J	0.63	0/1700	0.85	0/2320
3	R	0.66	0/71	0.87	0/94
3	S	0.65	0/84	0.76	0/112
3	T	0.52	0/86	0.73	0/115
3	V	0.61	0/78	0.79	0/104
All	All	0.61	0/14120	0.85	9/19206 (0.0%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	129	ALA	CA-C-N	6.67	126.63	119.76
2	I	129	ALA	C-N-CA	6.67	126.63	119.76
1	Q	218	GLU	N-CA-C	-6.20	105.81	113.19
1	A	63	VAL	N-CA-C	5.49	113.35	108.63
1	L	180	LEU	CA-CB-CG	5.45	135.36	116.30
2	B	99	ASP	CA-C-N	5.17	128.87	120.60
2	B	99	ASP	C-N-CA	5.17	128.87	120.60
2	H	219	SER	N-CA-C	5.08	116.82	111.28
1	P	180	LEU	CA-CB-CG	5.04	133.92	116.30

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	1713	0	1663	47	0
1	L	1702	0	1650	37	0
1	P	1722	0	1667	36	0
1	Q	1726	0	1673	36	0
2	B	1656	0	1605	18	0
2	H	1662	0	1607	31	0
2	I	1648	0	1601	27	0
2	J	1660	0	1608	28	0
3	R	71	0	67	2	0
3	S	84	0	81	4	0
3	T	86	0	87	2	0
3	V	78	0	76	3	0
4	A	6	0	8	2	0
4	J	6	0	8	3	0
4	P	6	0	8	7	0
4	Q	6	0	8	0	0
5	A	45	0	0	3	0
5	B	53	0	0	3	0
5	H	49	0	0	2	0
5	I	15	0	0	1	0
5	J	35	0	0	1	0
5	L	30	0	0	1	0
5	P	49	0	0	4	0
5	Q	35	0	0	2	0
5	S	1	0	0	0	0
All	All	14144	0	13417	253	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 9.

All (253) 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:45:PRO:HG2	1:L:170:GLU:HG2	1.27	1.13
1:A:6:GLN:HG2	1:A:23[A]:CYS:SG	1.93	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:2:ILE:HD11	1:L:98:LEU:HD12	1.41	1.01
2:B:57:THR:HG21	3:S:10:GLN:OE1	1.63	0.97
1:L:43:GLN:HE22	2:H:39:GLN:HE22	1.18	0.91
1:L:45:PRO:CG	1:L:170:GLU:HG2	2.04	0.87
1:Q:29:LEU:O	1:Q:97:THR:HG21	1.77	0.84
2:J:171:PRO:HA	4:J:401:GOL:H11	1.59	0.84
1:L:6:GLN:HG3	1:L:23[A]:CYS:SG	2.18	0.83
2:I:87:ARG:O	2:I:115:VAL:HG11	1.80	0.81
1:L:2:ILE:HD11	1:L:98:LEU:CD1	2.12	0.80
1:L:6:GLN:CG	1:L:23[A]:CYS:SG	2.69	0.80
1:P:6:GLN:HG3	1:P:23[A]:CYS:SG	2.23	0.79
2:B:54:LEU:HD11	3:S:15:ARG:HG3	1.65	0.78
2:H:83:MET:HE2	2:H:86:LEU:HD21	1.66	0.76
1:Q:43:GLN:HE22	2:J:39:GLN:HE22	1.33	0.76
1:P:43:GLN:HE22	2:I:39:GLN:HE22	1.33	0.75
1:P:45:PRO:HG3	1:P:170:GLU:HG2	1.69	0.74
1:A:110:GLU:HG3	1:A:178:TYR:OH	1.87	0.74
1:A:85:PRO:HA	1:A:111:ILE:HG21	1.69	0.74
2:I:204:HIS:HD2	2:I:207:SER:OG	1.69	0.73
1:L:88:PHE:CZ	1:L:170:GLU:HG3	2.24	0.72
2:I:12:VAL:HG21	2:I:18:LEU:HG	1.70	0.72
2:B:98:ARG:HD2	2:B:105:ASP:OD1	1.91	0.71
1:Q:130:LEU:O	1:Q:188:LYS:HD3	1.89	0.71
1:P:203:HIS:CD2	1:P:205:GLY:H	2.09	0.71
2:J:155:THR:HG22	5:J:227:HOH:O	1.90	0.71
1:A:42:GLN:HE22	4:A:404:GOL:H32	1.58	0.69
2:B:30:SER:O	2:B:53[A]:ASP:OD1	2.11	0.69
2:I:98:ARG:HD2	2:I:105:ASP:OD1	1.92	0.69
2:B:151:PRO:O	2:B:204:HIS:HE1	1.76	0.68
1:L:85:PRO:HA	1:L:111:ILE:HD12	1.76	0.68
2:H:151:PRO:O	2:H:204:HIS:HE1	1.77	0.68
1:A:2:ILE:HD11	1:A:98:LEU:HD13	1.74	0.67
1:A:203:HIS:CD2	1:A:205:GLY:H	2.12	0.67
1:P:169:THR:HG21	5:P:220:HOH:O	1.95	0.67
1:L:39:HIS:HD2	1:L:55:LYS:H	1.43	0.66
1:A:43:GLN:HE22	2:B:39:GLN:HE22	1.43	0.66
1:P:169:THR:HG22	1:P:170:GLU:O	1.95	0.66
1:L:6:GLN:HG2	1:L:23[A]:CYS:SG	2.35	0.65
1:P:203:HIS:HD2	1:P:205:GLY:H	1.43	0.65
2:H:71:SER:CB	4:P:402:GOL:H12	2.28	0.64
1:Q:216:ARG:HG2	1:Q:217:GLY:N	2.12	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:6:GLN:HG2	1:Q:23[A]:CYS:SG	2.38	0.64
2:J:204:HIS:HD2	2:J:207:SER:OG	1.81	0.64
1:A:117:ALA:HB1	1:A:206:LEU:HD13	1.80	0.63
1:L:88:PHE:HZ	1:L:170:GLU:HG3	1.63	0.63
2:B:83:MET:HE2	2:B:86:LEU:HD21	1.79	0.63
2:I:83:MET:HE2	2:I:86:LEU:HD21	1.81	0.63
2:J:172:ALA:H	4:J:401:GOL:C1	2.12	0.63
1:Q:203:HIS:CD2	1:Q:205:GLY:H	2.16	0.63
1:A:3:GLN:HG2	1:A:26:SER:HB3	1.80	0.62
1:P:37:TYR:HB2	1:P:97:THR:HG22	1.81	0.62
1:P:117:ALA:HB1	1:P:206:LEU:HD13	1.81	0.62
2:I:155:THR:HG22	5:I:233:HOH:O	1.99	0.62
3:T:13:PRO:HG2	3:T:16:VAL:HG23	1.81	0.62
1:A:29:LEU:O	1:A:97:THR:HG21	2.00	0.62
1:P:68:SER:HB2	4:P:402:GOL:H11	1.82	0.61
1:P:88:PHE:HZ	1:P:170:GLU:HG3	1.65	0.61
1:L:39:HIS:CD2	1:L:55:LYS:H	2.17	0.61
1:L:38:LEU:HG	1:L:39:HIS:N	2.12	0.61
2:J:151:PRO:O	2:J:204:HIS:HE1	1.84	0.60
1:Q:6:GLN:CG	1:Q:23[A]:CYS:SG	2.89	0.60
2:B:204:HIS:HD2	2:B:207:SER:OG	1.85	0.59
1:Q:203:HIS:HD2	1:Q:205:GLY:H	1.50	0.59
2:H:100:ASN:HB2	2:H:103:ALA:H	1.67	0.59
1:P:68:SER:HB2	4:P:402:GOL:C1	2.32	0.59
1:Q:38:LEU:HD22	1:Q:76:PHE:CG	2.37	0.59
2:J:64:VAL:HG13	2:J:68:PHE:HB2	1.83	0.59
1:P:88:PHE:CZ	1:P:170:GLU:HG3	2.38	0.58
1:Q:2:ILE:HD11	1:Q:98:LEU:HD13	1.85	0.58
1:P:6:GLN:NE2	1:P:91:TYR:O	2.36	0.58
5:P:297:HOH:O	2:I:133:LYS:HB2	2.03	0.58
1:P:196:VAL:HG22	1:P:215:ASN:ND2	2.18	0.58
2:H:204:HIS:HD2	2:H:207:SER:OG	1.87	0.58
2:B:1:GLU:O	2:B:26:GLY:HA3	2.04	0.58
2:B:32:TYR:O	2:B:72:ARG:NH2	2.38	0.57
2:I:98:ARG:CD	2:I:105:ASP:OD1	2.52	0.57
1:A:13:ALA:HB3	1:A:83:LEU:HD22	1.86	0.57
2:J:83:MET:HE2	2:J:86:LEU:HD21	1.86	0.57
2:H:71:SER:HB2	4:P:402:GOL:H12	1.87	0.57
1:A:39:HIS:CD2	1:A:55:LYS:H	2.23	0.56
1:A:74:THR:HG22	1:A:75:ASP:OD2	2.04	0.56
1:A:96:ASN:ND2	5:A:241:HOH:O	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:HIS:HD2	1:A:205:GLY:H	1.51	0.56
1:A:39:HIS:HD2	1:A:55:LYS:H	1.52	0.56
1:P:29:LEU:O	1:P:97:THR:HG21	2.05	0.56
1:A:85:PRO:HA	1:A:111:ILE:HD12	1.88	0.56
2:H:98:ARG:HD2	2:H:105:ASP:OD1	2.06	0.56
2:I:72:ARG:HD3	2:I:74:ASN:OD1	2.06	0.55
1:Q:98:LEU:HD23	3:V:7:ALA:HB1	1.89	0.55
2:J:4:LEU:HB3	2:J:96[A]:CYS:SG	2.46	0.55
1:A:45:PRO:CB	1:A:170:GLU:HG3	2.37	0.55
1:P:216:ARG:HG2	1:P:217:GLY:N	2.22	0.55
2:H:214:LYS:HE3	2:H:216:GLU:HG3	1.88	0.54
1:A:111:ILE:HG12	1:A:111:ILE:O	2.06	0.54
2:B:155:THR:HG22	5:B:227:HOH:O	2.08	0.54
2:J:53:ASP:HB2	2:J:101:TRP:HZ2	1.73	0.54
2:J:98:ARG:HD2	2:J:105:ASP:OD1	2.06	0.54
1:P:2:ILE:HD11	1:P:98:LEU:HD13	1.90	0.54
2:I:100:ASN:O	2:I:101:TRP:HB2	2.06	0.54
1:L:153:TRP:CD1	1:L:164:SER:HG	2.26	0.53
2:I:123:PRO:HB3	2:I:149:TYR:HB3	1.90	0.53
2:J:32:TYR:O	2:J:72:ARG:NH2	2.41	0.53
1:A:6:GLN:CG	1:A:23[A]:CYS:SG	2.84	0.53
1:P:45:PRO:CG	1:P:170:GLU:HG2	2.38	0.52
1:A:36:THR:HG21	5:A:243:HOH:O	2.08	0.52
1:Q:216:ARG:HG2	1:Q:217:GLY:H	1.74	0.52
1:Q:39:HIS:HE1	1:Q:96:ASN:HB2	1.73	0.52
1:L:33:ASN:HD22	1:L:33:ASN:C	2.17	0.52
2:H:3:GLN:HG2	2:H:25:SER:HB3	1.92	0.52
2:H:64:VAL:HG13	2:H:68:PHE:HB2	1.92	0.52
1:A:206:LEU:HG	1:A:210:VAL:HG23	1.92	0.52
2:H:12:VAL:HG21	2:H:18:LEU:HG	1.92	0.52
2:J:99:ASP:OD2	3:V:11:ARG:NH2	2.27	0.52
1:Q:37:TYR:HB2	1:Q:97:THR:HG22	1.92	0.51
2:H:83:MET:HE2	2:H:86:LEU:CD2	2.40	0.51
1:P:196:VAL:HG22	1:P:215:ASN:HD21	1.75	0.51
1:Q:111:ILE:HG12	1:Q:171:GLN:OE1	2.10	0.51
2:B:53[A]:ASP:HB2	2:B:101:TRP:HZ2	1.76	0.51
1:P:69:GLY:O	4:P:402:GOL:O2	2.26	0.51
1:A:74:THR:HG21	5:A:221:HOH:O	2.11	0.51
1:L:6:GLN:O	1:L:105:GLN:NE2	2.44	0.50
2:J:100:ASN:HB2	2:J:103:ALA:H	1.77	0.50
1:A:45:PRO:HB3	1:A:170:GLU:HG3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:151:PRO:O	2:I:204:HIS:HE1	1.94	0.50
1:L:95:GLN:C	1:L:96:ASN:HD22	2.19	0.50
1:A:39:HIS:HD2	1:A:54:TYR:HA	1.77	0.50
2:J:53:ASP:HB2	2:J:101:TRP:CZ2	2.47	0.50
1:L:13:ALA:HB3	1:L:83:LEU:HD22	1.94	0.50
1:P:169:THR:CG2	5:P:220:HOH:O	2.58	0.50
2:H:32:TYR:O	2:H:72:ARG:NH2	2.41	0.50
3:S:11:ARG:HA	3:S:11:ARG:NE	2.27	0.49
2:I:64:VAL:HG13	2:I:68:PHE:HB2	1.93	0.49
1:Q:134:THR:HB	5:Q:259:HOH:O	2.11	0.49
1:Q:180:LEU:HD12	1:Q:181:SER:N	2.27	0.49
1:Q:43:GLN:NE2	2:J:39:GLN:HE22	2.04	0.49
1:A:166:GLU:HB3	1:A:180:LEU:HD11	1.95	0.49
1:Q:37:TYR:HB2	1:Q:97:THR:CG2	2.43	0.48
1:P:191:TYR:CZ	1:P:216:ARG:HG3	2.49	0.48
1:P:39:HIS:HD2	1:P:55:LYS:H	1.61	0.48
2:J:109:GLN:CD	2:J:109:GLN:H	2.21	0.48
1:L:14:SER:O	1:L:17:ASP:HB2	2.13	0.48
1:A:72:SER:O	1:A:74:THR:N	2.46	0.48
1:A:38:LEU:HG	1:A:39:HIS:N	2.26	0.48
1:Q:111:ILE:HD12	1:Q:176:SER:OG	2.13	0.48
1:A:113:ARG:HG2	1:A:114:THR:N	2.29	0.48
1:P:59[A]:ARG:HD3	5:P:249:HOH:O	2.13	0.48
1:L:203:HIS:CD2	1:L:205:GLY:H	2.32	0.48
2:H:53[B]:ASP:CG	3:R:14:ASP:H	2.22	0.48
2:J:34:ILE:CG2	2:J:79:LEU:HD22	2.43	0.48
1:P:38:LEU:HD13	1:P:76:PHE:CD1	2.49	0.47
2:I:4:LEU:HB3	2:I:96:CYS:SG	2.54	0.47
1:Q:117:ALA:HB1	1:Q:206:LEU:HD13	1.97	0.47
1:Q:146:PRO:O	1:Q:203:HIS:HE1	1.97	0.47
2:H:53[B]:ASP:OD2	3:R:14:ASP:N	2.45	0.47
1:A:168:VAL:HG22	1:A:180:LEU:HD22	1.97	0.47
1:L:145:TYR:CG	1:L:146:PRO:HA	2.51	0.46
1:P:39:HIS:CD2	1:P:55:LYS:H	2.33	0.46
1:A:113:ARG:HG2	1:A:114:THR:H	1.80	0.46
1:Q:6:GLN:HG3	1:Q:23[A]:CYS:SG	2.56	0.46
2:H:6:GLU:HG3	2:H:96[A]:CYS:SG	2.55	0.46
1:Q:191:TYR:CZ	1:Q:216:ARG:HG3	2.51	0.46
1:A:3:GLN:HG2	1:A:26:SER:CB	2.46	0.46
1:L:168:VAL:HG22	1:L:180:LEU:HD22	1.97	0.46
1:Q:152[B]:GLN:OE1	1:Q:154:LYS:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:PRO:O	1:A:203:HIS:HE1	1.99	0.46
2:I:205:LYS:N	2:I:206:PRO:CD	2.79	0.46
1:P:71:GLY:HA3	1:P:76:PHE:HA	1.98	0.45
2:I:9:GLY:HA3	2:I:111:THR:OG1	2.16	0.45
1:A:147[B]:ARG:CZ	1:A:168:VAL:HG21	2.46	0.45
2:H:53[B]:ASP:OD1	2:H:53[B]:ASP:N	2.49	0.45
2:H:168:HIS:HB2	2:H:185:VAL:HG12	1.97	0.45
2:H:199:ILE:HG13	2:H:214:LYS:HA	1.98	0.45
2:B:62:ASP:O	2:B:65:THR:HG22	2.17	0.45
1:A:42:GLN:OE1	4:A:404:GOL:H11	2.17	0.45
2:H:68:PHE:HA	2:H:82:GLN:O	2.17	0.45
2:B:68:PHE:HA	2:B:82:GLN:O	2.16	0.45
2:H:72:ARG:HD3	2:H:74:ASN:OD1	2.17	0.44
2:B:132:SER:HB2	5:B:282:HOH:O	2.17	0.44
2:J:204:HIS:CD2	2:J:207:SER:OG	2.66	0.44
2:I:63:THR:HG23	2:I:64:VAL:HG23	2.00	0.44
3:T:11:ARG:NE	3:T:11:ARG:HA	2.32	0.44
1:A:39:HIS:CD2	1:A:54:TYR:HA	2.52	0.44
1:Q:37:TYR:HD1	1:Q:97:THR:HG22	1.83	0.44
1:L:60:PHE:CZ	2:H:105:ASP:HB2	2.53	0.44
2:I:88:ALA:HA	2:I:115:VAL:HG13	2.00	0.44
2:J:120:THR:HA	2:J:150:PHE:O	2.17	0.44
1:A:45:PRO:CG	1:A:170:GLU:HG3	2.48	0.44
2:J:83:MET:HE1	2:J:94:TYR:CZ	2.53	0.44
2:H:51:ILE:HD11	4:P:402:GOL:H2	1.99	0.44
1:A:43:GLN:NE2	2:B:39:GLN:HE22	2.11	0.43
2:I:68:PHE:HA	2:I:82:GLN:O	2.17	0.43
1:L:39:HIS:HD2	1:L:55:LYS:N	2.13	0.43
2:B:56:TYR:CD1	3:S:14:ASP:HB3	2.53	0.43
1:Q:49:PRO:HG2	2:J:107:TRP:CZ3	2.53	0.43
2:H:139:THR:HG22	5:H:283:HOH:O	2.18	0.43
2:I:100:ASN:HB2	2:I:103:ALA:H	1.84	0.43
1:Q:134:THR:HG22	1:Q:135:ALA:N	2.33	0.43
1:Q:130:LEU:HD12	1:Q:188:LYS:HB2	2.00	0.43
2:H:100:ASN:ND2	5:H:253:HOH:O	2.51	0.43
1:P:168:VAL:HG22	1:P:180:LEU:HD22	2.01	0.43
1:P:120:VAL:HA	1:P:140:LEU:O	2.17	0.43
1:L:156:ASP:O	1:L:157:ASN:HB2	2.19	0.43
1:Q:39:HIS:HD2	1:Q:55:LYS:H	1.67	0.43
2:H:143:GLY:HA2	2:H:158:TRP:CH2	2.54	0.42
2:J:130:PRO:HG3	2:J:142:LEU:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:38:LEU:HD22	1:L:76:PHE:CG	2.54	0.42
1:L:29:LEU:O	1:L:97:THR:HG21	2.19	0.42
1:P:52:LEU:HA	1:P:63:VAL:HG21	2.00	0.42
2:I:2:VAL:HG13	2:I:27:PHE:CD1	2.53	0.42
1:A:29:LEU:HD12	1:A:76:PHE:CE1	2.55	0.42
2:I:22:CYS:HB3	2:I:79:LEU:HB3	2.01	0.42
1:L:108:LYS:HG2	5:L:239:HOH:O	2.20	0.42
2:H:83:MET:HB3	2:H:86:LEU:HD21	2.02	0.42
1:L:146:PRO:O	1:L:203:HIS:HE1	2.02	0.42
1:A:113:ARG:HD3	1:A:114:THR:O	2.20	0.42
2:H:205:LYS:HE3	5:B:285:HOH:O	2.20	0.42
1:Q:10:SER:HA	1:Q:108:LYS:O	2.20	0.42
1:L:111:ILE:HG13	1:L:176:SER:OG	2.20	0.41
1:P:88:PHE:O	1:P:89:ALA:HB2	2.20	0.41
2:I:32:TYR:O	2:I:72:ARG:NH2	2.50	0.41
1:P:188:LYS:HE3	1:P:192[A]:GLU:OE2	2.20	0.41
1:P:43:GLN:NE2	2:I:39:GLN:HE22	2.07	0.41
1:Q:37:TYR:CE1	3:V:11:ARG:HG3	2.54	0.41
1:Q:25:SER:HB3	1:Q:27:GLN:O	2.20	0.41
1:L:53:ILE:HA	1:L:58:ASN:O	2.20	0.41
2:H:100:ASN:O	2:H:101:TRP:HB2	2.21	0.41
2:B:151:PRO:O	2:B:204:HIS:CE1	2.65	0.41
2:J:34:ILE:HG21	2:J:79:LEU:HD22	2.01	0.41
1:L:113:ARG:HG2	1:L:114:THR:N	2.36	0.41
1:A:168:VAL:HG22	1:A:180:LEU:HD13	2.03	0.41
1:A:63:VAL:HA	1:A:64:PRO:HD3	1.96	0.41
2:J:68:PHE:HA	2:J:82:GLN:O	2.21	0.41
1:L:2:ILE:CD1	1:L:98:LEU:HD12	2.30	0.41
2:H:188:VAL:HB	2:H:189:PRO:HD2	2.03	0.41
1:A:20:THR:HG23	1:A:77:THR:HG23	2.03	0.41
2:J:3:GLN:HG3	2:J:25:SER:HB3	2.03	0.41
1:A:150:LYS:HB3	1:A:202:THR:HB	2.01	0.41
2:I:216:GLU:OE2	2:I:217:PRO:HD2	2.20	0.41
1:L:105:GLN:H	1:L:105:GLN:HG3	1.60	0.40
1:A:2:ILE:CD1	1:A:98:LEU:HD13	2.49	0.40
2:I:52:SER:O	2:I:53:ASP:C	2.65	0.40
1:Q:25:SER:HB2	1:Q:74:THR:HA	2.03	0.40
1:P:39:HIS:HE1	1:P:96:ASN:HB2	1.85	0.40
1:Q:43:GLN:HE22	2:J:39:GLN:NE2	2.09	0.40
1:Q:166:GLU:HG3	5:Q:241:HOH:O	2.20	0.40
2:J:172:ALA:N	4:J:401:GOL:C1	2.83	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:31:HIS:HB3	1:L:33:ASN:ND2	2.36	0.40
1:P:68:SER:HB2	4:P:402:GOL:H32	2.03	0.40
1:L:85:PRO:O	1:L:111:ILE:HD12	2.21	0.40
1:A:12:SER:OG	1:A:110:GLU:OE2	2.40	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	219/219 (100%)	209 (95%)	9 (4%)	1 (0%)	24	37
1	L	218/219 (100%)	208 (95%)	10 (5%)	0	100	100
1	P	220/219 (100%)	210 (96%)	10 (4%)	0	100	100
1	Q	221/219 (101%)	214 (97%)	7 (3%)	0	100	100
2	B	219/226 (97%)	206 (94%)	12 (6%)	1 (0%)	24	37
2	H	220/226 (97%)	206 (94%)	11 (5%)	3 (1%)	9	13
2	I	218/226 (96%)	206 (94%)	10 (5%)	2 (1%)	14	22
2	J	220/226 (97%)	207 (94%)	10 (4%)	3 (1%)	9	13
3	R	7/35 (20%)	7 (100%)	0	0	100	100
3	S	9/35 (26%)	8 (89%)	1 (11%)	0	100	100
3	T	9/35 (26%)	9 (100%)	0	0	100	100
3	V	8/35 (23%)	7 (88%)	1 (12%)	0	100	100
All	All	1788/1920 (93%)	1697 (95%)	81 (4%)	10 (1%)	21	32

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	I	53	ASP
2	J	53	ASP
2	H	133	LYS
2	H	131	SER
1	A	73	GLY
2	H	138	GLY
2	B	133	LYS
2	I	133	LYS
2	J	131	SER
2	J	138	GLY

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	198/196 (101%)	173 (87%)	25 (13%)	4	6
1	L	197/196 (100%)	178 (90%)	19 (10%)	8	12
1	P	199/196 (102%)	180 (90%)	19 (10%)	8	13
1	Q	200/196 (102%)	183 (92%)	17 (8%)	10	16
2	B	184/189 (97%)	165 (90%)	19 (10%)	7	11
2	H	185/189 (98%)	171 (92%)	14 (8%)	12	21
2	I	183/189 (97%)	170 (93%)	13 (7%)	13	23
2	J	185/189 (98%)	169 (91%)	16 (9%)	10	16
3	R	7/27 (26%)	7 (100%)	0	100	100
3	S	9/27 (33%)	9 (100%)	0	100	100
3	T	9/27 (33%)	8 (89%)	1 (11%)	6	9
3	V	8/27 (30%)	7 (88%)	1 (12%)	4	6
All	All	1564/1648 (95%)	1420 (91%)	144 (9%)	8	14

All (144) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	1	ASP

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Mol	Chain	Res	Type
1	L	5	THR
1	L	33	ASN
1	L	36	THR
1	L	55	LYS
1	L	97	THR
1	L	105	GLN
1	L	111	ILE
1	L	130	LEU
1	L	137	VAL
1	L	140	LEU
1	L	161	SER
1	L	164	SER
1	L	180	LEU
1	L	186	LEU
1	L	190	ASP
1	L	206	LEU
1	L	208	SER
1	L	219	CYS
2	H	12	VAL
2	H	17	SER
2	H	20	LEU
2	H	34	ILE
2	H	43	LYS
2	H	64	VAL
2	H	98	ARG
2	H	116	SER
2	H	174	LEU
2	H	179	LEU
2	H	182	LEU
2	H	185	VAL
2	H	199	ILE
2	H	213	LYS
1	A	7	SER
1	A	12	SER
1	A	23[A]	CYS
1	A	23[B]	CYS
1	A	27	GLN
1	A	30	VAL
1	A	36	THR
1	A	38	LEU
1	A	63	VAL
1	A	74	THR

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Mol	Chain	Res	Type
1	A	97	THR
1	A	108	LYS
1	A	111	ILE
1	A	112	LYS
1	A	119	SER
1	A	128	GLU
1	A	130	LEU
1	A	137	VAL
1	A	140	LEU
1	A	159	LEU
1	A	180	LEU
1	A	186	LEU
1	A	206	LEU
1	A	208	SER
1	A	212	LYS
2	B	4	LEU
2	B	7	SER
2	B	12	VAL
2	B	43	LYS
2	B	57	THR
2	B	64	VAL
2	B	87	ARG
2	B	98	ARG
2	B	102	ASP
2	B	139	THR
2	B	142	LEU
2	B	147	LYS
2	B	155	THR
2	B	173	VAL
2	B	174	LEU
2	B	176	SER
2	B	179	LEU
2	B	185	VAL
2	B	219	SER
1	P	1	ASP
1	P	3	GLN
1	P	20	THR
1	P	23[A]	CYS
1	P	23[B]	CYS
1	P	36	THR
1	P	38	LEU
1	P	61	SER

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Mol	Chain	Res	Type
1	P	70	SER
1	P	74	THR
1	P	97	THR
1	P	98	LEU
1	P	130	LEU
1	P	140	LEU
1	P	159	LEU
1	P	180	LEU
1	P	193	LYS
1	P	206	LEU
1	P	219	CYS
2	I	12	VAL
2	I	17	SER
2	I	43	LYS
2	I	64	VAL
2	I	87	ARG
2	I	98	ARG
2	I	115	VAL
2	I	139	THR
2	I	154	VAL
2	I	155	THR
2	I	174	LEU
2	I	179	LEU
2	I	213	LYS
1	Q	36	THR
1	Q	70	SER
1	Q	74	THR
1	Q	94	SER
1	Q	96	ASN
1	Q	98	LEU
1	Q	111	ILE
1	Q	128	GLU
1	Q	130	LEU
1	Q	140	LEU
1	Q	159	LEU
1	Q	180	LEU
1	Q	186	LEU
1	Q	188	LYS
1	Q	206	LEU
1	Q	208	SER
1	Q	219	CYS
2	J	12	VAL

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Mol	Chain	Res	Type
2	J	20	LEU
2	J	25	SER
2	J	43	LYS
2	J	48	VAL
2	J	64	VAL
2	J	71	SER
2	J	98	ARG
2	J	139	THR
2	J	142	LEU
2	J	155	THR
2	J	174	LEU
2	J	179	LEU
2	J	185	VAL
2	J	199	ILE
2	J	205	LYS
3	T	9	SER
3	V	9	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	33	ASN
1	L	39	HIS
1	L	84	GLN
1	L	96	ASN
1	L	152	GLN
1	L	203	HIS
1	L	215	ASN
2	H	39	GLN
2	H	77	ASN
2	H	204	HIS
1	A	39	HIS
1	A	43	GLN
1	A	152	GLN
1	A	203	HIS
2	B	82	GLN
2	B	84	ASN
2	B	204	HIS
1	P	32	ASN
1	P	43	GLN
1	P	84	GLN
1	P	203	HIS

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Mol	Chain	Res	Type
1	P	215	ASN
2	I	204	HIS
1	Q	39	HIS
1	Q	43	GLN
1	Q	58	ASN
1	Q	84	GLN
1	Q	96	ASN
1	Q	203	HIS
1	Q	215	ASN
2	J	175	GLN
2	J	204	HIS

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](#)

4 ligands are modelled in this entry.

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	$\# Z > 2$
4	GOL	P	402	-	5,5,5	0.59	0	5,5,5	1.27	0
4	GOL	Q	403	-	5,5,5	0.43	0	5,5,5	0.22	0
4	GOL	A	404	-	5,5,5	0.50	0	5,5,5	0.48	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	J	401	-	5,5,5	0.38	0	5,5,5	0.65	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
4	GOL	P	402	-	-	0/4/4/4	-
4	GOL	Q	403	-	-	1/4/4/4	-
4	GOL	A	404	-	-	3/4/4/4	-
4	GOL	J	401	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	404	GOL	O1-C1-C2-C3
4	J	401	GOL	O2-C2-C3-O3
4	J	401	GOL	O1-C1-C2-C3
4	J	401	GOL	C1-C2-C3-O3
4	A	404	GOL	O1-C1-C2-O2
4	J	401	GOL	O1-C1-C2-O2
4	Q	403	GOL	O1-C1-C2-O2
4	A	404	GOL	C1-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	P	402	GOL	7	0
4	A	404	GOL	2	0
4	J	401	GOL	3	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	219/219 (100%)	0.44	9 (4%) 41 37	6, 39, 52, 67	6 (2%)
1	L	219/219 (100%)	0.49	6 (2%) 56 52	7, 39, 53, 67	4 (1%)
1	P	219/219 (100%)	0.79	12 (5%) 30 27	7, 39, 52, 66	7 (3%)
1	Q	219/219 (100%)	0.71	12 (5%) 30 27	6, 39, 53, 70	8 (3%)
2	B	220/226 (97%)	0.78	13 (5%) 28 24	13, 37, 53, 103	9 (4%)
2	H	220/226 (97%)	1.03	16 (7%) 21 17	12, 37, 55, 105	10 (4%)
2	I	220/226 (97%)	1.36	16 (7%) 21 17	6, 37, 54, 111	10 (4%)
2	J	220/226 (97%)	0.86	13 (5%) 28 24	13, 37, 52, 108	11 (5%)
3	R	9/35 (25%)	0.48	1 (11%) 10 7	37, 42, 54, 62	0
3	S	11/35 (31%)	1.01	2 (18%) 3 3	37, 42, 46, 80	1 (9%)
3	T	11/35 (31%)	0.87	1 (9%) 15 12	37, 40, 46, 53	0
3	V	10/35 (28%)	1.11	2 (20%) 3 2	37, 42, 51, 57	0
All	All	1797/1920 (93%)	0.81	103 (5%) 29 25	6, 38, 53, 111	66 (3%)

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	135	THR	37.9
2	I	134	SER	34.6
2	I	132	SER	34.5
1	P	218	GLU	33.2
2	I	133	LYS	31.9
2	I	136	SER	30.9
1	Q	218	GLU	28.1
2	I	131	SER	27.5
1	A	218	GLU	25.6
1	L	218	GLU	25.4
2	H	132	SER	25.1

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Mol	Chain	Res	Type	RSRZ
2	I	137	GLY	24.6
2	B	137	GLY	23.8
2	H	134	SER	23.7
2	J	135	THR	23.5
2	B	134	SER	23.3
2	H	131	SER	23.1
1	P	217	GLY	23.1
2	B	132	SER	22.9
2	B	136	SER	22.9
2	H	135	THR	22.0
2	B	135	THR	21.8
2	J	134	SER	21.4
2	J	136	SER	20.4
2	H	133	LYS	20.4
2	B	133	LYS	20.4
2	J	133	LYS	20.3
1	Q	219	CYS	18.9
2	H	136	SER	18.5
2	J	132	SER	17.6
2	J	137	GLY	17.0
1	L	217	GLY	16.5
1	P	219	CYS	15.9
1	Q	217	GLY	15.2
1	A	217	GLY	14.0
1	A	219	CYS	13.4
1	L	219	CYS	13.0
2	B	138	GLY	10.9
2	H	138	GLY	10.2
2	H	137	GLY	10.2
2	I	1	GLU	7.6
1	A	1	ASP	6.9
2	I	138	GLY	6.9
3	V	16	VAL	5.7
1	Q	24	ARG	5.6
2	J	1	GLU	5.4
3	S	16	VAL	5.0
2	I	219	SER	4.9
2	J	220	CYS	4.7
2	I	220	CYS	4.7
2	J	131	SER	4.6
1	P	1	ASP	4.6
3	S	15	ARG	4.4

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Mol	Chain	Res	Type	RSRZ
2	J	219	SER	4.4
1	Q	23[A]	CYS	4.2
2	I	53	ASP	4.1
2	J	138	GLY	4.1
2	H	220	CYS	3.8
1	Q	111	ILE	3.8
2	B	131	SER	3.8
1	L	23[A]	CYS	3.8
2	B	194	GLY	3.5
1	P	23[A]	CYS	3.5
1	Q	12	SER	3.3
2	B	220	CYS	3.2
2	H	219	SER	3.1
1	P	96	ASN	3.1
2	H	198	TYR	3.0
1	A	2	ILE	3.0
2	I	102	ASP	2.8
1	P	195	LYS	2.8
2	B	55	ALA	2.7
1	A	131	LYS	2.7
2	H	216	GLU	2.6
2	B	173	VAL	2.6
1	Q	74	THR	2.6
1	A	27	GLN	2.6
1	Q	186	LEU	2.5
1	Q	76	PHE	2.4
1	P	12	SER	2.4
2	I	194	GLY	2.4
1	L	96	ASN	2.4
2	H	1	GLU	2.4
1	P	135	ALA	2.3
2	J	194	GLY	2.3
1	A	23[A]	CYS	2.3
1	P	37	TYR	2.3
2	H	100	ASN	2.3
2	H	173	VAL	2.3
2	I	56	TYR	2.3
3	V	7	ALA	2.3
3	T	7	ALA	2.2
2	J	195	THR	2.2
3	R	15	ARG	2.2
1	L	134	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	96	ASN	2.2
1	Q	98	LEU	2.2
2	H	139	THR	2.1
1	Q	96	ASN	2.1
1	P	208	SER	2.1
1	P	186	LEU	2.1
2	B	87	ARG	2.1
2	I	29	PHE	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](#)

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(Å ²)	Q<0.9
4	GOL	A	404	6/6	0.79	0.17	60,62,62,63	0
4	GOL	Q	403	6/6	0.81	0.16	49,53,56,57	0
4	GOL	J	401	6/6	0.87	0.11	45,49,51,53	0
4	GOL	P	402	6/6	0.92	0.23	34,40,44,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.