



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:19 PM UTC

PDB ID : 7XY8 / pdb_00007xy8
Title : Crystal structure of antibody Fab fragment in complex with CD147(EMMPIRIN)
Authors : Nakamura, K.; Amano, M.; Yoneda, K.; Suzuki, M.; Fukuchi, K.
Deposited on : 2022-06-01
Resolution : 2.30 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:

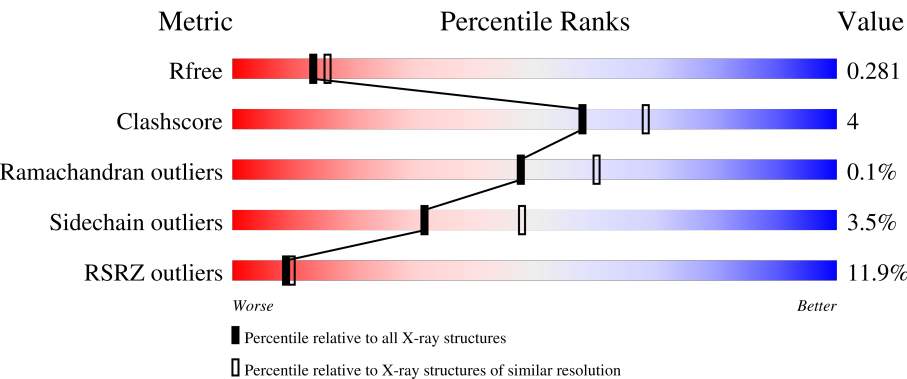
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	205	49% 75%12%•12%
1	B	205	8% 41%7%51%
2	H	236	3% 80%10%10%
2	I	236	82%10%•8%
3	L	214	4% 87%11%•

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	M	214	% 84% 15%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8889 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 Isoform 2 of Basigin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	181	Total	C	N	O	S	0	0	0
			1389	862	238	281	8			
1	B	100	Total	C	N	O	S	0	0	0
			768	471	135	157	5			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP P35613
A	2	GLY	-	expression tag	UNP P35613
A	3	SER	-	expression tag	UNP P35613
A	4	SER	-	expression tag	UNP P35613
A	5	HIS	-	expression tag	UNP P35613
A	6	HIS	-	expression tag	UNP P35613
A	7	HIS	-	expression tag	UNP P35613
A	8	HIS	-	expression tag	UNP P35613
A	9	HIS	-	expression tag	UNP P35613
A	10	HIS	-	expression tag	UNP P35613
A	11	SER	-	expression tag	UNP P35613
A	12	SER	-	expression tag	UNP P35613
A	13	GLY	-	expression tag	UNP P35613
A	14	GLU	-	expression tag	UNP P35613
A	15	ASN	-	expression tag	UNP P35613
A	16	LEU	-	expression tag	UNP P35613
A	17	TYR	-	expression tag	UNP P35613
A	18	PHE	-	expression tag	UNP P35613
A	19	GLN	-	expression tag	UNP P35613
A	20	GLY	-	expression tag	UNP P35613
A	21	SER	-	expression tag	UNP P35613
B	1	MET	-	initiating methionine	UNP P35613
B	2	GLY	-	expression tag	UNP P35613
B	3	SER	-	expression tag	UNP P35613
B	4	SER	-	expression tag	UNP P35613

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	5	HIS	-	expression tag	UNP P35613
B	6	HIS	-	expression tag	UNP P35613
B	7	HIS	-	expression tag	UNP P35613
B	8	HIS	-	expression tag	UNP P35613
B	9	HIS	-	expression tag	UNP P35613
B	10	HIS	-	expression tag	UNP P35613
B	11	SER	-	expression tag	UNP P35613
B	12	SER	-	expression tag	UNP P35613
B	13	GLY	-	expression tag	UNP P35613
B	14	GLU	-	expression tag	UNP P35613
B	15	ASN	-	expression tag	UNP P35613
B	16	LEU	-	expression tag	UNP P35613
B	17	TYR	-	expression tag	UNP P35613
B	18	PHE	-	expression tag	UNP P35613
B	19	GLN	-	expression tag	UNP P35613
B	20	GLY	-	expression tag	UNP P35613
B	21	SER	-	expression tag	UNP P35613

- Molecule 2 is a protein called heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	212	Total	C	N	O	S	0	0	0
			1604	1008	272	315	9			
2	I	218	Total	C	N	O	S	0	0	0
			1649	1032	281	327	9			

- Molecule 3 is a protein called light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	214	Total	C	N	O	S	0	0	0
			1646	1026	277	337	6			
3	M	214	Total	C	N	O	S	0	0	0
			1646	1026	277	337	6			

- Molecule 4 is water.

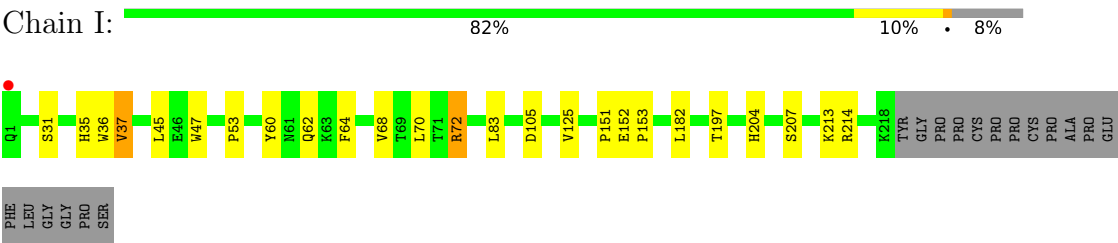
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	5	Total	O	0	0
			5	5		
4	B	8	Total	O	0	0
			8	8		

Continued on next page...

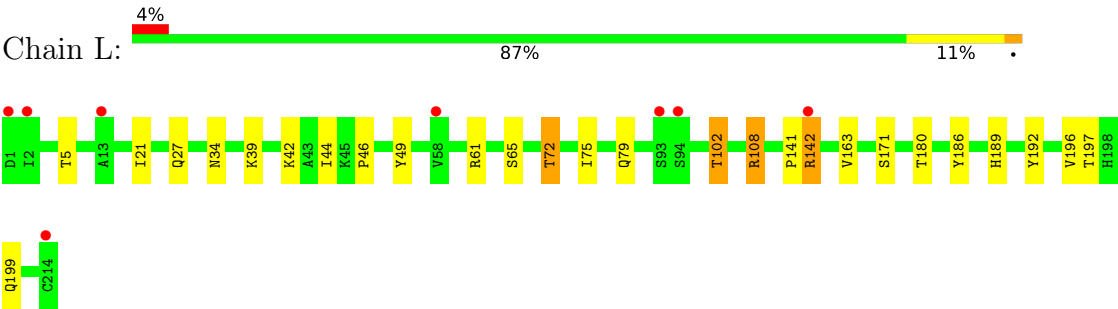
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	16	Total 16	O 16	0	0
4	I	74	Total 74	O 74	0	0
4	L	27	Total 27	O 27	0	0
4	M	57	Total 57	O 57	0	0

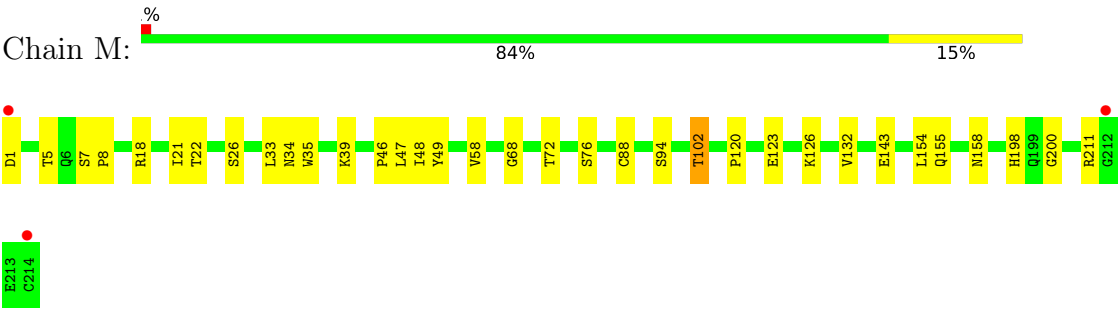
● Molecule 2: heavy chain



● Molecule 3: light chain



● Molecule 3: light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.96Å 93.37Å 98.31Å 90.00° 90.89° 90.00°	Depositor
Resolution (Å)	35.91 – 2.30 35.91 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.1 (35.91-2.30) 98.1 (35.91-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.236 , 0.281 0.236 , 0.281	Depositor DCC
R_{free} test set	2502 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	33.0	Xtriage
Anisotropy	0.528	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 25.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.043 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8889	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.32% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.58	0/1416	0.86	2/1918 (0.1%)
1	B	0.54	0/782	0.84	0/1056
2	H	0.51	0/1642	0.84	0/2234
2	I	0.53	0/1688	0.83	0/2297
3	L	0.53	0/1681	0.86	1/2281 (0.0%)
3	M	0.53	0/1681	0.85	0/2281
All	All	0.54	0/8890	0.85	3/12067 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	103	GLY	CA-C-N	6.90	124.63	119.66
1	A	103	GLY	C-N-CA	6.90	124.63	119.66
3	L	189	HIS	N-CA-C	5.04	117.91	110.46

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	A	1389	0	1343	11	0
1	B	768	0	739	7	0
2	H	1604	0	1562	13	0
2	I	1649	0	1604	16	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	L	1646	0	1591	13	0
3	M	1646	0	1591	22	0
4	A	5	0	0	0	0
4	B	8	0	0	0	0
4	H	16	0	0	0	0
4	I	74	0	0	0	0
4	L	27	0	0	1	0
4	M	57	0	0	6	0
All	All	8889	0	8430	77	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 4.

All (77) 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)
3:M:5:THR:HB	4:M:312:HOH:O	1.64	0.97
2:I:35:HIS:HD2	2:I:47:TRP:HE1	1.32	0.74
3:M:198:HIS:CD2	3:M:200:GLY:H	2.07	0.72
3:M:5:THR:CB	4:M:312:HOH:O	2.28	0.69
3:M:21:ILE:HG23	3:M:102:THR:HG21	1.79	0.65
3:M:155:GLN:HE21	3:M:158:ASN:HD21	1.44	0.65
3:M:5:THR:CG2	4:M:312:HOH:O	2.44	0.63
3:M:26:SER:HB2	4:M:351:HOH:O	1.98	0.62
2:H:35:HIS:HD2	2:H:47:TRP:HE1	1.46	0.62
3:M:8:PRO:O	3:M:102:THR:HB	2.01	0.61
3:L:61:ARG:HH21	3:L:79:GLN:HE21	1.49	0.60
2:H:105:ASP:HA	3:L:46:PRO:HG3	1.84	0.59
2:I:151:PRO:O	2:I:204:HIS:HE1	1.86	0.59
1:B:142:ILE:HB	1:B:182:GLN:HB2	1.85	0.57
3:M:5:THR:HG22	4:M:312:HOH:O	2.04	0.57
2:H:11:VAL:HG21	2:H:151:PRO:HG3	1.87	0.56
1:B:175:ASN:HD21	1:B:178:ALA:H	1.52	0.55
3:L:142:ARG:NH2	3:L:163:VAL:HG21	2.22	0.54
3:M:120:PRO:HD3	3:M:132:VAL:HG22	1.90	0.54
3:M:22:THR:HG22	3:M:72:THR:HG22	1.89	0.53
1:A:159:PHE:HB2	1:A:170:HIS:HB2	1.91	0.53
2:H:39:GLN:HB2	2:H:45:LEU:HD13	1.89	0.53
3:M:18:ARG:HG3	3:M:76:SER:HA	1.89	0.53
1:B:136:ASP:HB3	1:B:188:THR:HB	1.91	0.52
3:M:33:LEU:HD11	3:M:88:CYS:HB2	1.91	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:VAL:HA	1:A:132:PRO:C	2.35	0.52
1:A:29:THR:HB	1:A:40:THR:HB	1.91	0.51
2:I:152:GLU:HG3	2:I:153:PRO:HA	1.93	0.51
3:M:21:ILE:CG2	3:M:102:THR:HG21	2.39	0.51
3:M:198:HIS:HD2	3:M:200:GLY:H	1.58	0.50
3:L:34:ASN:HD22	3:L:49:TYR:HA	1.76	0.50
2:I:125:VAL:O	2:I:213:LYS:HE3	2.11	0.50
1:B:131:VAL:HA	1:B:132:PRO:C	2.37	0.49
2:H:35:HIS:CD2	2:H:47:TRP:HE1	2.28	0.49
1:B:122:ALA:HB2	1:B:174:LEU:HD11	1.95	0.48
2:I:35:HIS:CD2	2:I:47:TRP:HE1	2.21	0.48
1:A:54:ARG:HB3	1:A:88:VAL:HB	1.95	0.48
2:I:204:HIS:HD2	2:I:207:SER:OG	1.97	0.47
3:M:68:GLY:HA2	4:M:344:HOH:O	2.13	0.47
2:I:105:ASP:HA	3:M:46:PRO:HG3	1.96	0.47
3:L:141:PRO:HD2	3:L:199:GLN:HG3	1.97	0.47
3:M:35:TRP:HB2	3:M:48:ILE:HB	1.97	0.47
2:H:151:PRO:O	2:H:204:HIS:HE1	1.98	0.47
1:A:26:VAL:HG13	1:A:97:ALA:HB2	1.98	0.46
1:A:81:GLN:HA	1:A:101:LEU:HD13	1.96	0.46
3:L:108:ARG:HG3	3:L:171:SER:HB2	1.96	0.46
1:B:110:VAL:HB	2:I:31:SER:HB2	1.98	0.45
1:A:122:ALA:HB3	1:A:171:ILE:HB	1.98	0.45
3:M:34:ASN:HD22	3:M:49:TYR:HA	1.82	0.45
2:H:123:PRO:HB3	2:H:149:TYR:HB3	1.99	0.45
1:A:177:GLU:HA	1:A:178:ALA:HA	1.75	0.45
3:M:123:GLU:HA	3:M:126:LYS:HE3	1.99	0.45
3:L:21:ILE:CG2	3:L:102:THR:HG21	2.47	0.44
2:I:37:VAL:CG1	2:I:45:LEU:HD22	2.48	0.44
1:B:129:GLU:HG2	2:I:35:HIS:CE1	2.53	0.43
2:H:35:HIS:CD2	2:H:50:TRP:HB3	2.52	0.43
2:I:64:PHE:HB3	2:I:68:VAL:HG23	2.00	0.43
2:H:48:MET:HE2	2:H:68:VAL:HG21	1.99	0.43
1:A:27:PHE:HB3	1:A:42:SER:HB2	2.00	0.43
1:A:137:TRP:CD2	1:A:167:SER:HB2	2.54	0.43
2:H:154:VAL:HG12	2:H:204:HIS:HB2	2.00	0.43
1:A:92:GLU:HA	1:A:93:PRO:HA	1.77	0.43
2:I:60:TYR:HE1	2:I:70:LEU:HG	1.84	0.43
3:L:199:GLN:H	3:L:199:GLN:HG2	1.65	0.43
2:H:67:ARG:NH2	2:H:90:ASP:OD2	2.52	0.42
2:H:107:TRP:CG	3:L:44:ILE:HB	2.54	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:M:47:LEU:HA	3:M:58:VAL:HG21	2.02	0.42
2:I:197:THR:HG22	2:I:214:ARG:HH21	1.84	0.41
3:L:180:THR:HB	4:L:325:HOH:O	2.21	0.41
3:L:186:TYR:HA	3:L:192:TYR:OH	2.21	0.41
3:L:65:SER:HB3	3:L:72:THR:HG23	2.03	0.41
2:I:53:PRO:HA	2:I:72:ARG:HG2	2.02	0.41
2:H:190:SER:HA	2:H:193:LEU:HD13	2.02	0.41
2:I:36:TRP:CD1	2:I:70:LEU:HD22	2.56	0.41
3:M:7:SER:HA	3:M:8:PRO:HA	1.86	0.41
3:L:39:LYS:HB2	3:L:42:LYS:HB2	2.04	0.40
2:I:152:GLU:HG3	2:I:153:PRO:CA	2.51	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	179/205 (87%)	164 (92%)	14 (8%)	1 (1%)	21	27
1	B	98/205 (48%)	93 (95%)	5 (5%)	0	100	100
2	H	208/236 (88%)	205 (99%)	3 (1%)	0	100	100
2	I	216/236 (92%)	214 (99%)	2 (1%)	0	100	100
3	L	212/214 (99%)	202 (95%)	10 (5%)	0	100	100
3	M	212/214 (99%)	205 (97%)	7 (3%)	0	100	100
All	All	1125/1310 (86%)	1083 (96%)	41 (4%)	1 (0%)	48	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	190	SER

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	156/176 (89%)	151 (97%)	5 (3%)	34	51
1	B	87/176 (49%)	84 (97%)	3 (3%)	32	49
2	H	179/199 (90%)	174 (97%)	5 (3%)	38	56
2	I	185/199 (93%)	180 (97%)	5 (3%)	39	58
3	L	189/189 (100%)	180 (95%)	9 (5%)	23	35
3	M	189/189 (100%)	182 (96%)	7 (4%)	30	45
All	All	985/1128 (87%)	951 (96%)	34 (4%)	32	48

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	VAL
1	A	49	GLU
1	A	101	LEU
1	A	174	LEU
1	A	195	GLN
1	B	116	ILE
1	B	168	GLU
1	B	195	GLN
2	H	63	LYS
2	H	72	ARG
2	H	83	LEU
2	H	155	THR
2	H	182	LEU
2	I	37	VAL
2	I	62	GLN
2	I	72	ARG
2	I	83	LEU
2	I	182	LEU
3	L	5	THR
3	L	27	GLN
3	L	72	THR
3	L	75	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	L	102	THR
3	L	108	ARG
3	L	142	ARG
3	L	196	VAL
3	L	197	THR
3	M	1	ASP
3	M	39	LYS
3	M	94	SER
3	M	102	THR
3	M	143	GLU
3	M	154	LEU
3	M	211	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	GLN
1	A	81	GLN
1	A	117	ASN
1	A	152	ASN
1	B	117	ASN
1	B	175	ASN
2	H	35	HIS
2	H	62	GLN
2	H	204	HIS
2	I	1	GLN
2	I	35	HIS
2	I	39	GLN
2	I	62	GLN
2	I	175	GLN
2	I	204	HIS
3	L	27	GLN
3	L	34	ASN
3	L	79	GLN
3	M	3	GLN
3	M	34	ASN
3	M	38	GLN
3	M	55	GLN
3	M	155	GLN
3	M	198	HIS
3	M	210	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	181/205 (88%)	2.30	101 (55%) 0 0	43, 73, 126, 141	0
1	B	100/205 (48%)	1.15	16 (16%) 5 5	32, 51, 90, 102	0
2	H	212/236 (89%)	0.49	7 (3%) 49 51	22, 37, 58, 82	0
2	I	218/236 (92%)	-0.00	1 (0%) 87 87	17, 25, 42, 62	0
3	L	214/214 (100%)	0.53	8 (3%) 45 48	21, 37, 60, 93	0
3	M	214/214 (100%)	0.17	3 (1%) 73 75	17, 31, 45, 80	0
All	All	1139/1310 (86%)	0.69	136 (11%) 9 10	17, 37, 95, 141	0

All (136) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	30	VAL	6.8
1	A	23	ALA	6.0
1	A	33	LEU	5.6
1	A	67	LEU	5.5
1	A	39	LEU	5.4
1	A	37	ILE	5.2
1	A	38	LEU	5.0
3	M	214	CYS	4.8
1	A	47	ALA	4.7
3	M	212	GLY	4.6
1	A	178	ALA	4.4
1	A	27	PHE	4.4
1	A	42	SER	4.3
1	A	25	THR	4.3
1	A	28	THR	4.3
1	A	153	GLY	4.3
1	A	35	SER	4.1
1	A	48	THR	4.0
3	L	214	CYS	4.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	68	PRO	4.0
1	A	66	ALA	3.9
1	A	61	VAL	3.9
1	A	82	TRP	3.9
1	A	62	LEU	3.8
1	A	60	VAL	3.7
1	A	29	THR	3.6
1	A	89	PHE	3.5
1	A	43	LEU	3.5
1	B	116	ILE	3.5
1	A	76	VAL	3.5
1	A	170	HIS	3.5
1	A	74	PHE	3.4
2	H	130	PRO	3.3
1	A	58	GLY	3.3
1	A	88	VAL	3.3
1	A	149	ALA	3.3
1	A	99	ILE	3.3
2	H	131	CYS	3.2
1	A	72	THR	3.2
1	A	131	VAL	3.2
1	A	45	ASP	3.2
1	A	36	LYS	3.2
1	A	26	VAL	3.2
1	A	93	PRO	3.1
1	A	71	LYS	3.1
1	A	145	SER	3.1
1	A	24	GLY	3.1
1	A	87	CYS	3.1
1	A	51	THR	3.0
1	A	110	VAL	3.0
1	A	160	VAL	3.0
1	A	183	TYR	3.0
1	A	174	LEU	3.0
1	B	153	GLY	3.0
1	A	191	LYS	2.9
2	H	76	ILE	2.9
1	A	44	ASN	2.9
3	L	13	ALA	2.8
1	B	160	VAL	2.8
1	A	143	THR	2.8
1	A	63	LYS	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	50	VAL	2.8
1	A	137	TRP	2.8
1	A	79	ASP	2.8
1	A	102	HIS	2.8
1	A	141	LYS	2.8
1	A	139	TRP	2.8
1	A	158	PHE	2.8
1	A	41	CYS	2.7
1	A	34	GLY	2.7
1	A	78	SER	2.7
1	A	159	PHE	2.7
1	A	101	LEU	2.7
1	A	150	LEU	2.7
1	A	140	TYR	2.7
1	A	83	GLY	2.7
3	L	93	SER	2.6
1	A	53	HIS	2.6
1	A	31	GLU	2.6
1	B	118	GLU	2.6
1	A	192	GLY	2.6
1	B	159	PHE	2.6
1	B	143	THR	2.6
1	A	81	GLN	2.5
2	I	1	GLN	2.5
1	B	115	HIS	2.5
1	A	109	ALA	2.5
1	A	148	LYS	2.5
1	A	103	GLY	2.5
1	B	142	ILE	2.5
1	A	90	LEU	2.5
1	A	75	LYS	2.5
1	A	97	ALA	2.5
1	B	154	SER	2.4
1	A	146	GLU	2.4
2	H	42	GLY	2.4
1	A	151	MET	2.4
3	L	2	ILE	2.4
1	A	190	SER	2.4
1	A	59	GLY	2.3
1	B	144	ASP	2.3
1	A	85	TYR	2.3
2	H	29	PHE	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	H	54	GLY	2.3
1	B	158	PHE	2.3
3	L	58	VAL	2.3
1	A	32	ASP	2.2
1	A	154	SER	2.2
1	A	171	ILE	2.2
1	A	176	MET	2.2
3	L	1	ASP	2.2
1	A	49	GLU	2.2
1	B	178	ALA	2.2
1	A	152	ASN	2.1
1	A	55	TRP	2.1
1	A	133	PRO	2.1
2	H	53	PRO	2.1
1	A	69	GLY	2.1
3	L	94	SER	2.1
1	A	65	ASP	2.1
3	M	1	ASP	2.1
1	A	187	GLY	2.1
1	B	103	GLY	2.1
1	A	135	THR	2.1
1	A	94	MET	2.1
1	A	175	ASN	2.1
1	A	182	GLN	2.1
1	B	175	ASN	2.1
1	B	146	GLU	2.0
1	A	115	HIS	2.0
1	A	169	LEU	2.0
1	A	142	ILE	2.0
1	B	110	VAL	2.0
3	L	142	ARG	2.0
1	A	177	GLU	2.0
1	A	124	LEU	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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.