



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 16, 2022 – 08:18 PM JST

PDB ID : 7XY8
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	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

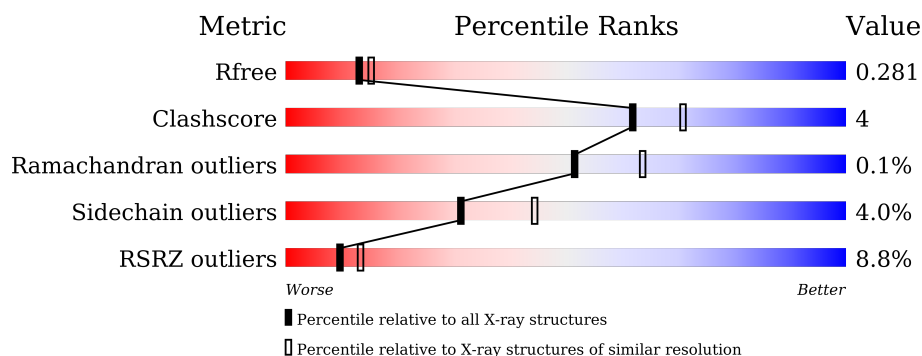
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (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	<div> <div>38%</div> <div> <div>74%</div> <div>14%</div> <div>•</div> <div>12%</div> </div> </div>
1	B	205	<div> <div>5%</div> <div> <div>41%</div> <div>7%</div> <div>51%</div> </div> </div>
2	H	236	<div> <div>2%</div> <div> <div>80%</div> <div>10%</div> <div>10%</div> </div> </div>
2	I	236	<div> <div>82%</div> <div> <div>10%</div> <div>•</div> <div>8%</div> </div> </div>
3	L	214	<div> <div>2%</div> <div> <div>88%</div> <div>10%</div> <div>•</div> </div> </div>
3	M	214	<div> <div>%</div> <div> <div>83%</div> <div>16%</div> </div> </div>

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

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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		

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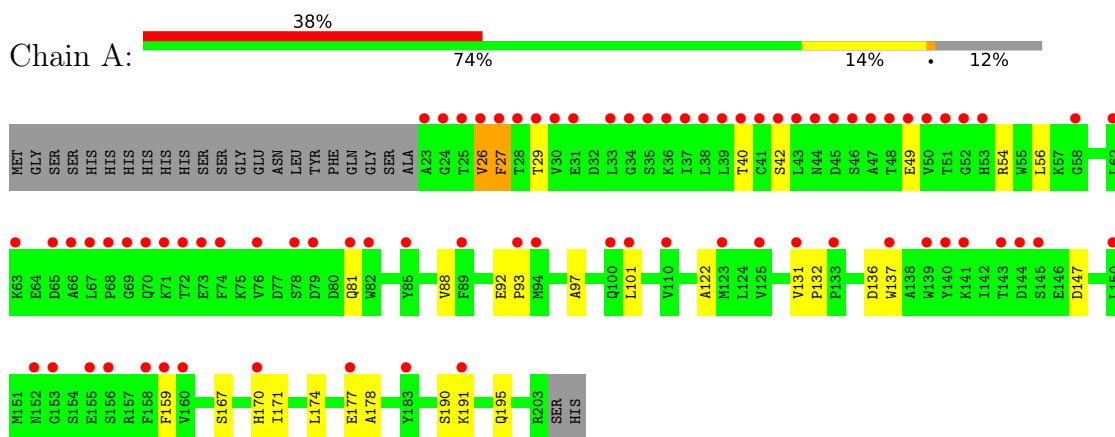
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	16	Total 16	O 16	0	0
4	I	74	Total 74	O 74	0	0
4	L	27	Total 27	O 27	0	0
4	M	57	Total 57	O 57	0	0

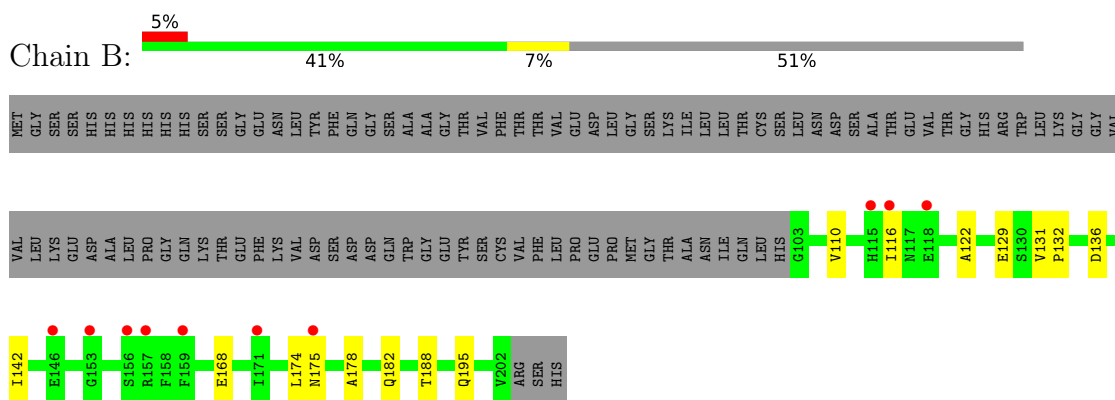
3 Residue-property plots [i](#)

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.

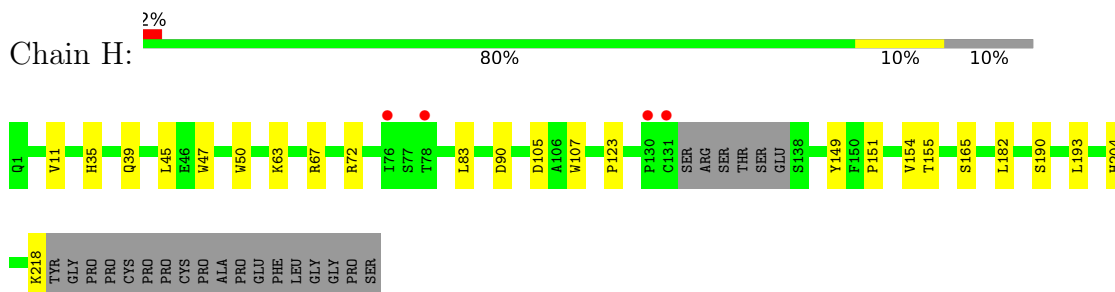
- Molecule 1: Isoform 2 of Basigin



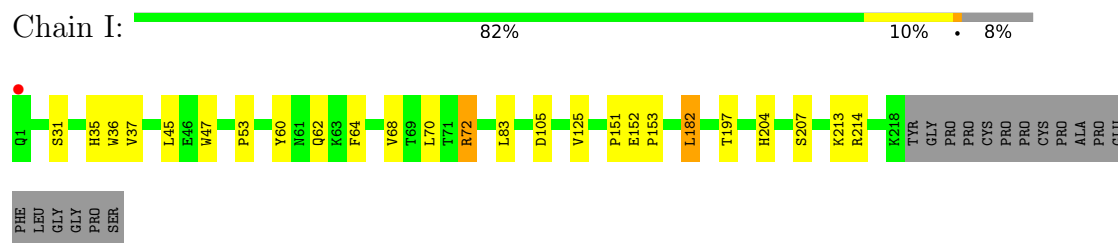
- Molecule 1: Isoform 2 of Basigin



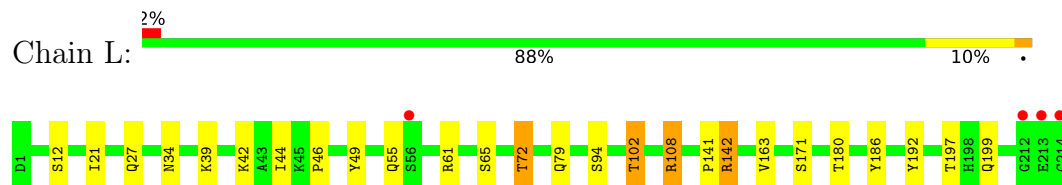
- Molecule 2: heavy chain



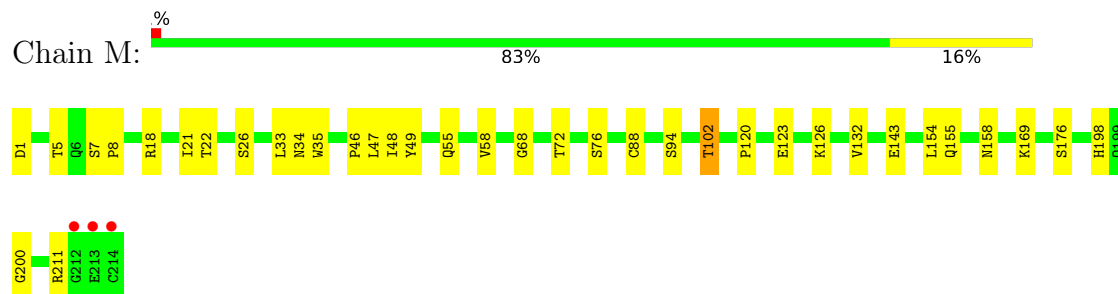
- Molecule 2: heavy chain



- 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 (4.88%)	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 , 29.5	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

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	0.41	0/1416	0.63	1/1918 (0.1%)
1	B	0.40	0/782	0.61	0/1056
2	H	0.41	0/1642	0.67	1/2234 (0.0%)
2	I	0.43	0/1688	0.67	1/2297 (0.0%)
3	L	0.41	0/1681	0.67	0/2281
3	M	0.43	0/1681	0.67	0/2281
All	All	0.42	0/8890	0.66	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(°)
2	I	182	LEU	CA-CB-CG	6.33	129.86	115.30
2	H	182	LEU	CA-CB-CG	5.82	128.68	115.30
1	A	56	LEU	CA-CB-CG	5.49	127.93	115.30

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	1389	0	1343	11	0
1	B	768	0	739	7	0
2	H	1604	0	1562	12	0
2	I	1649	0	1604	16	0

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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	76	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 (76) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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
1:A:131:VAL:HA	1:A:132:PRO:C	2.35	0.47
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
1:B:131:VAL:HA	1:B:132:PRO:C	2.37	0.44
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
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
3:M:47:LEU:HA	3:M:58:VAL:HG21	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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%)	25	31
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%)	51	64

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%)	148 (95%)	8 (5%)	24	33
1	B	87/176 (49%)	84 (97%)	3 (3%)	37	51
2	H	179/199 (90%)	173 (97%)	6 (3%)	37	51
2	I	185/199 (93%)	181 (98%)	4 (2%)	52	69
3	L	189/189 (100%)	180 (95%)	9 (5%)	25	36
3	M	189/189 (100%)	180 (95%)	9 (5%)	25	36
All	All	985/1128 (87%)	946 (96%)	39 (4%)	31	44

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	VAL
1	A	27	PHE
1	A	49	GLU
1	A	136	ASP
1	A	147	ASP
1	A	174	LEU
1	A	191	LYS
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	165	SER
2	H	218	LYS
2	I	62	GLN
2	I	72	ARG
2	I	83	LEU
2	I	182	LEU
3	L	12	SER

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Mol	Chain	Res	Type
3	L	27	GLN
3	L	55	GLN
3	L	72	THR
3	L	94	SER
3	L	102	THR
3	L	108	ARG
3	L	142	ARG
3	L	197	THR
3	M	1	ASP
3	M	55	GLN
3	M	94	SER
3	M	102	THR
3	M	143	GLU
3	M	154	LEU
3	M	169	LYS
3	M	176	SER
3	M	211	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	117	ASN
1	A	164	GLN
1	B	175	ASN
1	B	182	GLN
2	H	35	HIS
2	H	62	GLN
2	H	204	HIS
2	I	35	HIS
2	I	175	GLN
2	I	204	HIS
3	L	34	ASN
3	L	53	ASN
3	L	55	GLN
3	L	79	GLN
3	M	34	ASN
3	M	55	GLN
3	M	155	GLN
3	M	198	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 monosaccharides 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%)	2.16	78 (43%) 0 0	43, 73, 126, 141	0
1	B	100/205 (48%)	0.68	10 (10%) 7 10	32, 51, 90, 102	0
2	H	212/236 (89%)	0.14	4 (1%) 66 73	21, 37, 58, 82	0
2	I	218/236 (92%)	-0.02	1 (0%) 91 94	17, 25, 42, 62	0
3	L	214/214 (100%)	0.18	4 (1%) 66 73	21, 37, 60, 93	0
3	M	214/214 (100%)	0.02	3 (1%) 75 80	17, 31, 45, 80	0
All	All	1139/1310 (86%)	0.46	100 (8%) 10 13	17, 37, 95, 141	0

All (100) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	30	VAL	12.1
1	A	66	ALA	9.0
1	A	71	LYS	8.3
3	M	212	GLY	7.3
1	A	37	ILE	7.2
1	A	72	THR	7.2
1	A	24	GLY	6.8
1	A	73	GLU	6.1
1	A	145	SER	6.0
1	A	67	LEU	5.8
3	L	214	CYS	5.8
1	A	48	THR	5.6
1	A	27	PHE	5.5
1	A	51	THR	5.3
1	A	49	GLU	5.3
1	A	42	SER	5.1
1	A	63	LYS	4.9
1	A	38	LEU	4.8
1	A	46	SER	4.7

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Mol	Chain	Res	Type	RSRZ
1	A	69	GLY	4.6
1	A	25	THR	4.6
1	A	44	ASN	4.6
1	A	45	ASP	4.5
1	B	153	GLY	4.4
1	A	33	LEU	4.3
1	A	43	LEU	4.3
1	A	28	THR	4.3
1	A	82	TRP	4.3
1	A	40	THR	4.3
1	A	89	PHE	4.3
1	A	79	ASP	4.3
1	A	50	VAL	4.2
1	A	53	HIS	4.1
1	B	116	ILE	4.1
1	A	65	ASP	4.0
1	A	177	GLU	4.0
1	A	35	SER	3.9
1	A	170	HIS	3.8
1	A	153	GLY	3.7
1	A	68	PRO	3.7
1	A	41	CYS	3.7
1	A	62	LEU	3.7
1	A	39	LEU	3.7
3	M	214	CYS	3.6
1	A	47	ALA	3.6
1	A	140	TYR	3.6
2	H	76	ILE	3.4
1	A	152	ASN	3.4
1	B	146	GLU	3.4
1	A	23	ALA	3.3
1	A	159	PHE	3.3
1	A	160	VAL	3.3
1	A	70	GLN	3.2
2	H	131	CYS	3.2
1	A	101	LEU	3.1
1	A	131	VAL	3.0
1	A	125	VAL	3.0
1	A	36	LYS	2.9
1	A	26	VAL	2.9
1	A	85	TYR	2.8
1	A	144	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
2	H	130	PRO	2.8
1	B	159	PHE	2.8
1	A	31	GLU	2.8
1	A	76	VAL	2.7
1	A	78	SER	2.6
1	A	156	SER	2.6
1	A	29	THR	2.6
1	A	155	GLU	2.6
1	A	191	LYS	2.5
1	A	74	PHE	2.5
3	L	212	GLY	2.5
1	B	157	ARG	2.4
2	I	1	GLN	2.4
1	A	143	THR	2.4
1	B	118	GLU	2.4
1	A	81	GLN	2.4
1	A	123	MET	2.4
1	A	94	MET	2.3
1	B	115	HIS	2.3
1	A	100	GLN	2.3
1	A	150	LEU	2.3
1	B	156	SER	2.3
1	B	175	ASN	2.3
1	A	58	GLY	2.2
1	B	171	ILE	2.2
1	A	34	GLY	2.2
1	A	141	LYS	2.2
1	A	133	PRO	2.2
3	L	213	GLU	2.2
1	A	139	TRP	2.2
1	A	52	GLY	2.1
1	A	158	PHE	2.1
1	A	93	PRO	2.1
2	H	78	THR	2.1
1	A	110	VAL	2.1
1	A	183	TYR	2.1
1	A	137	TRP	2.1
3	L	56	SER	2.0
3	M	213	GLU	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 monosaccharides 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.