Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2024 – 01:17 PM EDT

PDB ID	:	1IFH
Title	:	A DETAILED ANALYSIS OF THE FREE AND BOUND CONFORMATION OF AN ANTIBODY: X-RAY STRUCTURES OF ANTI-PEPTIDE FAB 17(SLASH)9 AND THREE DIFFERENT FAB-PEPTIDE COMPLEXES
Authors	:	Schulze-Gahmen, U.; Wilson, I.A.
Deposited on	:	1993-05-06
Resolution	:	2.80 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:

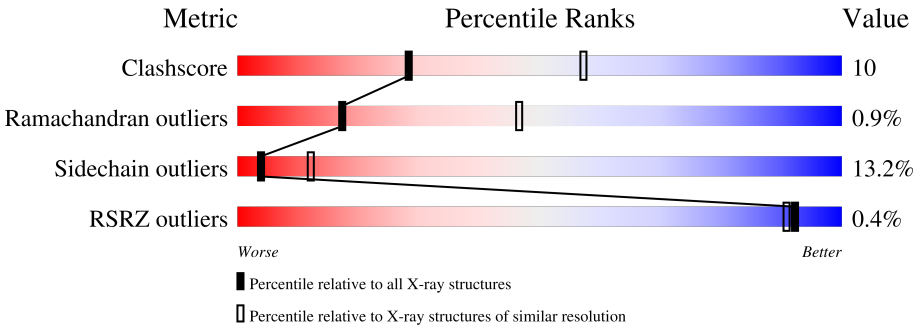
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.80 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	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 <=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	L	218	67% 26% 5% •
2	H	220	% 61% 32% 5% •
3	P	8	38% 50% 12%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3391 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 IGG2A-KAPPA 17/9 FAB (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	218	Total	C	N	O	S	0	0	1
			1680	1047	280	346	7			

- Molecule 2 is a protein called IGG2A-KAPPA 17/9 FAB (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	220	Total	C	N	O	S	0	0	1
			1653	1041	273	331	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	227	PRO	-	insertion	GB 533229

- Molecule 3 is a protein called INFLUENZA HEMAGGLUTININ HA1 (STRAIN X47) (RESIDUES 101-107).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	8	Total	C	N	O	0	0	0
			57	35	7	15			

- Molecule 4 is water.

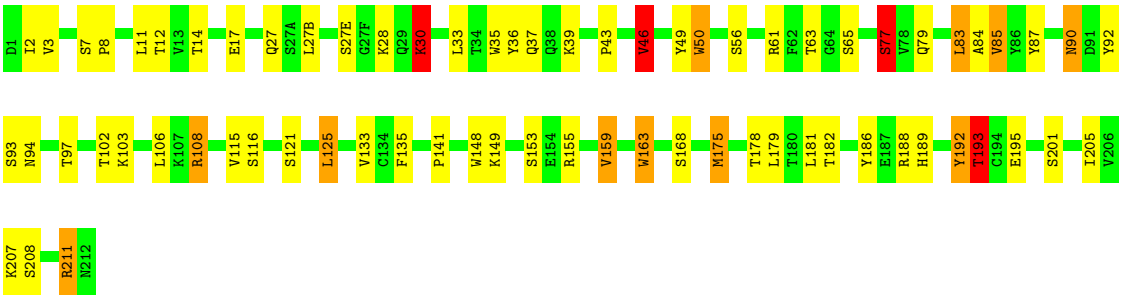
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	1	Total	O	0	0
			1	1		

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.

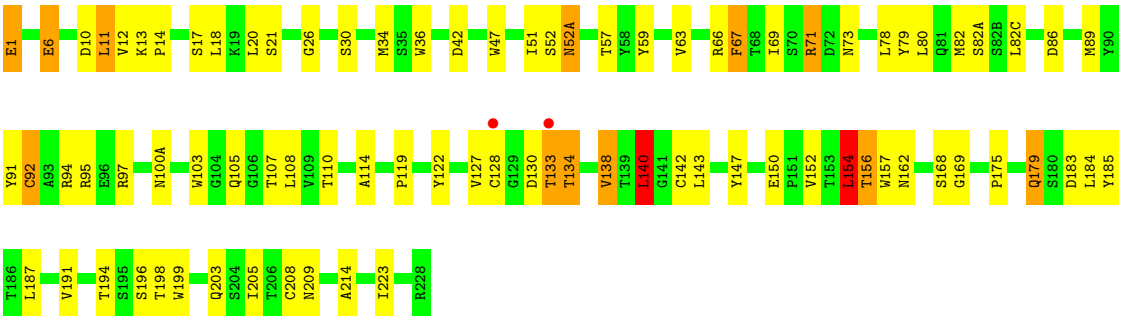
• Molecule 1: IGG2A-KAPPA 17/9 FAB (LIGHT CHAIN)

Chain L:



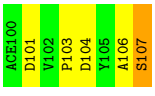
• Molecule 2: IGG2A-KAPPA 17/9 FAB (HEAVY CHAIN)

Chain H:



• Molecule 3: INFLUENZA HEMAGGLUTININ HA1 (STRAIN X47) (RESIDUES 101-107)

Chain P:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	73.10Å 103.90Å 67.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.80 44.61 – 2.67	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.80) 78.7 (44.61-2.67)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.67 (at 2.69Å)	Xtriage
Refinement program	X-PLOR 3.0	Depositor
R, R_{free}	0.170 , (Not available) 0.174 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	24.6	Xtriage
Anisotropy	0.201	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 102.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	3391	wwPDB-VP
Average B, all atoms (Å^2)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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)

Bond lengths and bond angles in the following residue types are not validated in this section: ACE

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	0.95	2/1718 (0.1%)	1.74	25/2336 (1.1%)
2	H	0.92	0/1694	1.78	40/2310 (1.7%)
3	P	1.01	1/56 (1.8%)	2.01	3/76 (3.9%)
All	All	0.94	3/3468 (0.1%)	1.76	68/4722 (1.4%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	107	SER	C-OXT	5.94	1.34	1.23
1	L	193	THR	CA-CB	5.46	1.67	1.53
1	L	50	TRP	CG-CD2	-5.39	1.34	1.43

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	59	TYR	CB-CG-CD2	-11.53	114.08	121.00
2	H	47	TRP	CD1-CG-CD2	10.13	114.41	106.30
2	H	157	TRP	CD1-CG-CD2	9.94	114.25	106.30
2	H	71	ARG	NE-CZ-NH1	9.48	125.04	120.30
1	L	148	TRP	CD1-CG-CD2	9.18	113.64	106.30
2	H	94	ARG	NE-CZ-NH1	8.64	124.62	120.30
2	H	66	ARG	NE-CZ-NH1	8.03	124.32	120.30
1	L	50	TRP	CD1-CG-CD2	8.03	112.72	106.30
1	L	148	TRP	CE2-CD2-CG	-7.99	100.91	107.30
2	H	47	TRP	CE2-CD2-CG	-7.91	100.97	107.30
1	L	61	ARG	NE-CZ-NH1	7.88	124.24	120.30
2	H	36	TRP	CD1-CG-CD2	7.64	112.41	106.30
1	L	188	ARG	NE-CZ-NH2	-7.55	116.52	120.30
2	H	199	TRP	CD1-CG-CD2	7.54	112.33	106.30
2	H	103	TRP	CE2-CD2-CG	-7.48	101.31	107.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	L	50	TRP	CE2-CD2-CG	-7.46	101.34	107.30
2	H	157	TRP	CE2-CD2-CG	-7.45	101.34	107.30
1	L	35	TRP	CD1-CG-CD2	7.42	112.24	106.30
2	H	140	LEU	CA-CB-CG	7.30	132.08	115.30
2	H	36	TRP	CE2-CD2-CG	-7.26	101.49	107.30
2	H	103	TRP	CD1-CG-CD2	7.17	112.04	106.30
1	L	192	TYR	CB-CG-CD2	-7.16	116.71	121.00
3	P	101	ASP	CA-C-N	-7.10	101.58	117.20
2	H	36	TRP	CG-CD2-CE3	6.77	139.99	133.90
1	L	35	TRP	CE2-CD2-CG	-6.76	101.89	107.30
1	L	163	TRP	CG-CD2-CE3	6.69	139.92	133.90
1	L	163	TRP	CE2-CD2-CG	-6.65	101.98	107.30
2	H	199	TRP	CE2-CD2-CG	-6.62	102.00	107.30
2	H	36	TRP	CB-CG-CD1	-6.61	118.41	127.00
2	H	107	THR	CA-CB-CG2	6.52	121.52	112.40
1	L	87	TYR	CB-CG-CD1	-6.42	117.14	121.00
2	H	157	TRP	CG-CD1-NE1	-6.40	103.70	110.10
1	L	46	VAL	CG1-CB-CG2	-6.31	100.81	110.90
2	H	71	ARG	NE-CZ-NH2	-6.21	117.19	120.30
2	H	34	MET	CG-SD-CE	-6.10	90.44	100.20
2	H	103	TRP	CG-CD2-CE3	6.08	139.37	133.90
1	L	85	VAL	CG1-CB-CG2	-6.05	101.23	110.90
1	L	77	SER	N-CA-C	-5.91	95.03	111.00
2	H	47	TRP	CG-CD1-NE1	-5.80	104.30	110.10
1	L	108	ARG	NE-CZ-NH1	5.78	123.19	120.30
2	H	185	TYR	CB-CG-CD1	-5.78	117.53	121.00
1	L	182	THR	CA-CB-CG2	-5.76	104.33	112.40
2	H	63	VAL	N-CA-CB	-5.75	98.85	111.50
2	H	79	TYR	CB-CG-CD2	-5.68	117.59	121.00
2	H	95	ARG	NE-CZ-NH1	5.64	123.12	120.30
2	H	154	LEU	CA-CB-CG	5.55	128.07	115.30
1	L	148	TRP	CG-CD1-NE1	-5.54	104.56	110.10
2	H	86	ASP	CB-CG-OD2	-5.48	113.37	118.30
2	H	103	TRP	CB-CG-CD1	-5.41	119.97	127.00
1	L	163	TRP	CB-CG-CD1	-5.38	120.01	127.00
2	H	66	ARG	NE-CZ-NH2	-5.37	117.61	120.30
1	L	30	LYS	CB-CA-C	-5.36	99.67	110.40
2	H	1	GLU	N-CA-C	-5.31	96.66	111.00
2	H	36	TRP	CG-CD1-NE1	-5.31	104.79	110.10
1	L	163	TRP	CD1-CG-CD2	5.25	110.50	106.30
1	L	108	ARG	NE-CZ-NH2	-5.25	117.67	120.30
2	H	6	GLU	CA-CB-CG	5.23	124.91	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	138	VAL	N-CA-C	-5.23	96.89	111.00
2	H	138	VAL	CG1-CB-CG2	-5.22	102.55	110.90
1	L	211	ARG	N-CA-C	5.21	125.06	111.00
2	H	156	THR	CA-CB-CG2	-5.19	105.13	112.40
3	P	101	ASP	CB-CG-OD2	-5.19	113.63	118.30
2	H	82(A)	SER	CA-C-N	-5.13	105.91	117.20
2	H	86	ASP	CB-CG-OD1	5.11	122.90	118.30
1	L	211	ARG	NE-CZ-NH1	5.07	122.83	120.30
1	L	92	TYR	CB-CG-CD2	-5.02	117.99	121.00
2	H	94	ARG	NE-CZ-NH2	-5.02	117.79	120.30
3	P	101	ASP	CB-CG-OD1	5.00	122.80	118.30

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	1680	0	1612	33	0
2	H	1653	0	1603	33	0
3	P	57	0	46	2	0
4	H	1	0	0	1	0
All	All	3391	0	3261	65	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 10.

All (65) 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)
2:H:51:ILE:HD13	2:H:71:ARG:HG2	1.74	0.70
1:L:2:ILE:HG21	1:L:27(B):LEU:HD21	1.75	0.69
2:H:133:THR:O	2:H:138:VAL:HG12	1.94	0.67
1:L:8:PRO:HG2	1:L:11:LEU:HD13	1.78	0.66
1:L:28:LYS:HE2	1:L:30:LYS:HE2	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:51:ILE:HG13	2:H:57:THR:HG22	1.80	0.64
2:H:119:PRO:HB3	2:H:147:TYR:HB3	1.79	0.64
3:P:104:ASP:O	3:P:107:SER:HB3	1.99	0.63
2:H:71:ARG:HD3	2:H:73:ASN:OD1	2.02	0.60
2:H:82:MET:HB3	2:H:82(C):LEU:HD21	1.85	0.59
1:L:36:TYR:CE1	1:L:46:VAL:HG23	2.39	0.58
1:L:193:THR:HB	1:L:208:SER:OG	2.02	0.58
1:L:163:TRP:CD1	1:L:175:MET:HG3	2.38	0.57
1:L:49:TYR:CD2	2:H:100(A):ASN:HB2	2.39	0.57
1:L:14:THR:O	1:L:17:GLU:HG2	2.05	0.56
2:H:127:VAL:HG13	2:H:128:CYS:H	1.72	0.55
2:H:67:PHE:HD2	2:H:80:LEU:HD21	1.70	0.55
2:H:12:VAL:HG11	2:H:82(C):LEU:HD12	1.89	0.55
1:L:43:PRO:HB3	2:H:91:TYR:CE2	2.42	0.54
1:L:115:VAL:HA	1:L:135:PHE:O	2.07	0.54
1:L:121:SER:O	1:L:125:LEU:HB2	2.08	0.53
2:H:122:TYR:HB2	2:H:143:LEU:HB3	1.90	0.53
2:H:30:SER:O	2:H:52(A):ASN:HB2	2.08	0.53
1:L:2:ILE:HG12	1:L:27:GLN:HB2	1.89	0.53
2:H:114:ALA:HB2	2:H:183:ASP:HB3	1.91	0.52
2:H:133:THR:HG23	2:H:134:THR:OG1	2.10	0.52
1:L:83:LEU:HD11	1:L:106:LEU:HB2	1.93	0.51
1:L:149:LYS:HA	1:L:153:SER:O	2.10	0.51
2:H:14:PRO:HA	2:H:82(C):LEU:O	2.10	0.51
1:L:90:ASN:HB3	1:L:97:THR:OG1	2.11	0.51
1:L:94:ASN:OD1	3:P:106:ALA:HB3	2.12	0.50
2:H:162:ASN:OD1	2:H:205:ILE:HA	2.11	0.50
2:H:11:LEU:HD22	2:H:110:THR:HB	1.95	0.49
1:L:85:VAL:HG22	1:L:103:LYS:HG3	1.95	0.49
1:L:85:VAL:HG22	1:L:103:LYS:CG	2.42	0.49
1:L:77:SER:HB3	1:L:79:GLN:HE21	1.78	0.49
2:H:1:GLU:O	2:H:26:GLY:HA3	2.13	0.48
1:L:36:TYR:HE1	1:L:46:VAL:HG23	1.78	0.48
1:L:133:VAL:HG22	1:L:178:THR:HG23	1.96	0.48
2:H:69:ILE:HD11	2:H:78:LEU:HD11	1.94	0.48
1:L:149:LYS:HB2	1:L:193:THR:HG23	1.96	0.47
1:L:159:VAL:HG12	1:L:179:LEU:HG	1.95	0.47
1:L:30:LYS:HE3	1:L:50:TRP:CE2	2.49	0.47
1:L:103:LYS:HB3	1:L:103:LYS:HE2	1.65	0.47
2:H:154:LEU:HD23	2:H:187:LEU:HD21	1.96	0.47
2:H:179:GLN:HG3	2:H:184:LEU:O	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:155:ARG:HG2	1:L:179:LEU:HD21	1.96	0.46
1:L:155:ARG:HA	1:L:155:ARG:HD2	1.68	0.46
2:H:67:PHE:CD2	2:H:80:LEU:HD21	2.49	0.46
2:H:169:GLY:O	2:H:191:VAL:HA	2.16	0.45
1:L:189:HIS:HB2	1:L:192:TYR:OH	2.17	0.45
2:H:20:LEU:HG	2:H:82:MET:CE	2.48	0.43
2:H:12:VAL:HG22	2:H:18:LEU:HD13	2.01	0.42
1:L:115:VAL:HG12	1:L:207:LYS:HG3	2.02	0.42
1:L:39:LYS:HD3	1:L:84:ALA:HB2	2.01	0.42
2:H:52:SER:HB2	4:H:229:HOH:O	2.19	0.42
1:L:8:PRO:O	1:L:102:THR:HG23	2.20	0.42
2:H:18:LEU:HB3	2:H:82:MET:HE3	2.01	0.42
2:H:6:GLU:HG3	2:H:92:CYS:SG	2.60	0.41
2:H:198:THR:O	2:H:203:GLN:HB2	2.21	0.41
1:L:186:TYR:CZ	1:L:211:ARG:HD3	2.56	0.41
2:H:20:LEU:HG	2:H:82:MET:HE2	2.02	0.41
1:L:195:GLU:HA	1:L:205:ILE:O	2.21	0.40
2:H:147:TYR:HE2	2:H:150:GLU:HG2	1.86	0.40
2:H:140:LEU:HB3	2:H:223:ILE:HG21	2.04	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	L	216/218 (99%)	209 (97%)	7 (3%)	0	100 100
2	H	218/220 (99%)	198 (91%)	17 (8%)	3 (1%)	11 34
3	P	6/8 (75%)	4 (67%)	1 (17%)	1 (17%)	0 0
All	All	440/446 (99%)	411 (93%)	25 (6%)	4 (1%)	17 46

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	133	THR
2	H	214	ALA
2	H	175	PRO
3	P	103	PRO

5.3.2 Protein sidechains (i)

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	194/195 (100%)	169 (87%)	25 (13%)	4 13
2	H	186/187 (100%)	160 (86%)	26 (14%)	3 11
3	P	6/6 (100%)	6 (100%)	0	100 100
All	All	386/388 (100%)	335 (87%)	51 (13%)	4 12

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	3	VAL
1	L	7	SER
1	L	12	THR
1	L	27(E)	SER
1	L	30	LYS
1	L	33	LEU
1	L	37	GLN
1	L	46	VAL
1	L	56	SER
1	L	63	THR
1	L	65	SER
1	L	77	SER
1	L	83	LEU
1	L	90	ASN
1	L	93	SER
1	L	108	ARG
1	L	116	SER
1	L	125	LEU
1	L	141	PRO

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Mol	Chain	Res	Type
1	L	159	VAL
1	L	168	SER
1	L	175	MET
1	L	181	LEU
1	L	193	THR
1	L	201	SER
2	H	10	ASP
2	H	11	LEU
2	H	13	LYS
2	H	17	SER
2	H	21	SER
2	H	42	ASP
2	H	52(A)	ASN
2	H	67	PHE
2	H	89	MET
2	H	92	CYS
2	H	97	ARG
2	H	105	GLN
2	H	108	LEU
2	H	130	ASP
2	H	134	THR
2	H	140	LEU
2	H	142	CYS
2	H	152	VAL
2	H	154	LEU
2	H	156	THR
2	H	168	SER
2	H	179	GLN
2	H	194	THR
2	H	196	SER
2	H	208	CYS
2	H	209	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	79	GLN
1	L	137	ASN
1	L	161	ASN
1	L	189	HIS
2	H	52(A)	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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, 95^{th} 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	L	218/218 (100%)	-0.74	0 100 100	2, 9, 35, 45	0
2	H	220/220 (100%)	-0.62	2 (0%) 84 80	2, 12, 49, 80	0
3	P	7/8 (87%)	-0.75	0 100 100	4, 7, 12, 24	0
All	All	445/446 (99%)	-0.68	2 (0%) 92 91	2, 10, 40, 80	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	133	THR	3.0
2	H	128	CYS	2.3

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no monosaccharides 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.