



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 05:07 pm BST

PDB ID : 1ADQ
Title : CRYSTAL STRUCTURE OF A HUMAN IGM RHEUMATOID FACTOR
FAB IN COMPLEX WITH ITS AUTOANTIGEN IGG FC
Authors : Corper, A.L.; Taussig, M.J.; Sutton, B.J.
Deposited on : 1997-02-18
Resolution : 3.15 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 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	:	2.11
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)	:	2.11

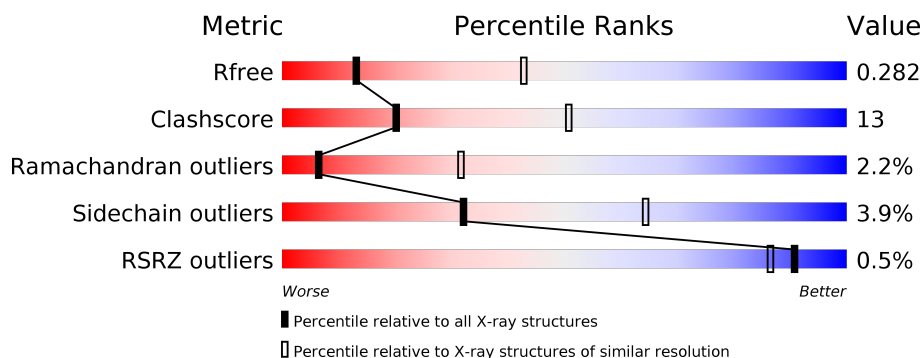
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 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	1665 (3.20-3.12)
Clashscore	141614	1804 (3.20-3.12)
Ramachandran outliers	138981	1770 (3.20-3.12)
Sidechain outliers	138945	1769 (3.20-3.12)
RSRZ outliers	127900	1616 (3.20-3.12)

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 on 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	206	
2	L	213	
3	H	225	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4977 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 IGG4 REA FC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	206	Total	C	N	O	S	381	0	0
			1653	1048	279	319	7			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	272	GLN	GLU	CONFLICT	UNP P01861
A	283	GLN	GLU	CONFLICT	UNP P01861
A	294	GLN	GLU	CONFLICT	UNP P01861
A	312	ASN	ASP	CONFLICT	UNP P01861
A	315	ASP	ASN	CONFLICT	UNP P01861

- Molecule 2 is a protein called IGM-LAMBDA RF-AN FAB (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	213	Total	C	N	O	S	193	0	0
			1588	991	267	325	5			

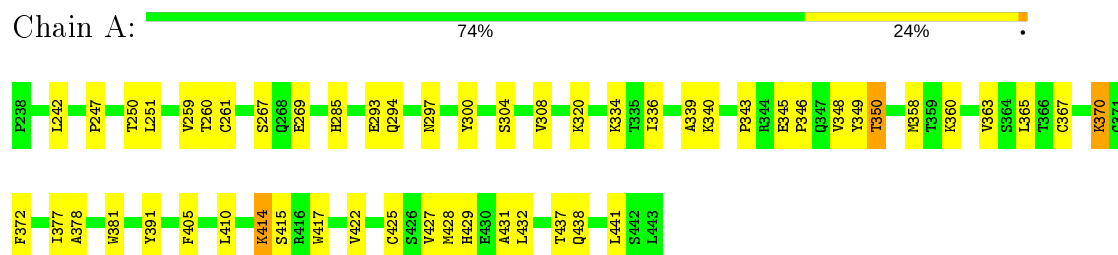
- Molecule 3 is a protein called IGM-LAMBDA RF-AN FAB (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	225	Total	C	N	O	S	481	0	0
			1736	1098	296	334	8			

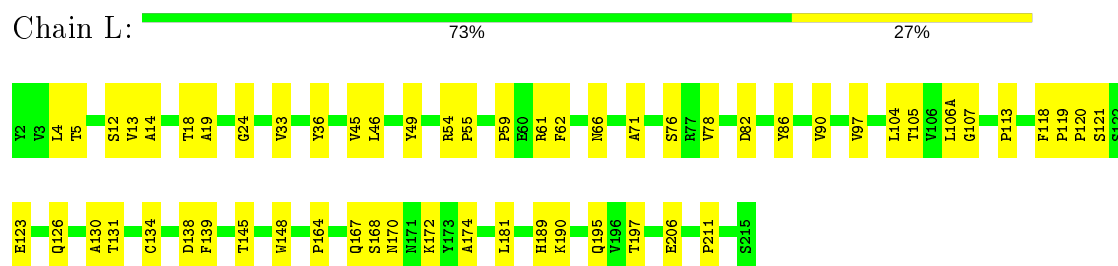
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

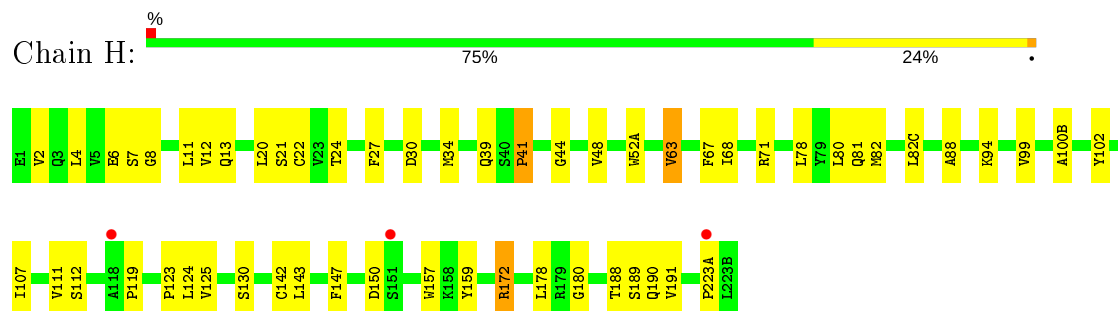
- Molecule 1: IGG4 REA FC



- Molecule 2: IGM-LAMBDA RF-AN FAB (LIGHT CHAIN)



- Molecule 3: IGM-LAMBDA RF-AN FAB (HEAVY CHAIN)



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	160.35Å 81.95Å 64.24Å 90.00° 98.27° 90.00°	Depositor
Resolution (Å)	8.00 – 3.15 8.00 – 3.15	Depositor EDS
% Data completeness (in resolution range)	86.0 (8.00-3.15) 86.0 (8.00-3.15)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.01 (at 3.14Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.225 , 0.289 0.220 , 0.282	Depositor DCC
R_{free} test set	756 reflections (6.52%)	wwPDB-VP
Wilson B-factor (Å ²)	54.1	Xtriage
Anisotropy	0.385	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 66.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	4977	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.43% 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	0.31	0/1697	0.56	0/2308
2	L	0.28	0/1629	0.53	0/2228
3	H	0.31	0/1778	0.56	0/2416
All	All	0.30	0/5104	0.55	0/6952

There are no bond length outliers.

There are no bond angle outliers.

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	1653	0	1615	33	0
2	L	1588	0	1523	33	0
3	H	1736	0	1692	31	0
All	All	4977	0	4830	95	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 13.

All (95) 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:H:119:PRO:HB3	3:H:147:PHE:HB3	1.46	0.97

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:8:GLY:HA3	3:H:20:LEU:HD23	1.67	0.74
1:A:259:VAL:HG23	1:A:308:VAL:HG21	1.70	0.73
3:H:82:MET:HB3	3:H:82(C):LEU:HD21	1.71	0.72
1:A:346:PRO:HB3	1:A:372:PHE:HB3	1.70	0.72
2:L:18:THR:HG22	2:L:76:SER:HA	1.71	0.72
3:H:34:MET:HB3	3:H:78:LEU:HD22	1.71	0.71
3:H:41:PRO:HD3	3:H:88:ALA:HA	1.72	0.70
1:A:348:VAL:HG21	1:A:427:VAL:HG11	1.79	0.62
3:H:30:ASP:O	3:H:52(A):TRP:HB2	1.99	0.62
1:A:242:LEU:HD23	1:A:336:ILE:HG23	1.82	0.60
2:L:13:VAL:HG21	2:L:19:ALA:HB2	1.83	0.60
1:A:432:LEU:HD13	1:A:437:THR:HG22	1.84	0.59
1:A:350:THR:HB	1:A:441:LEU:HG	1.84	0.59
2:L:121:SER:HB2	3:H:123:PRO:O	2.01	0.59
1:A:427:VAL:HG13	1:A:437:THR:HG23	1.85	0.56
2:L:46:LEU:O	2:L:55:PRO:HG3	2.06	0.55
2:L:138:ASP:HA	2:L:172:LYS:HB3	1.89	0.54
3:H:188:THR:HG22	3:H:189:SER:H	1.73	0.54
1:A:259:VAL:HG13	1:A:336:ILE:HD11	1.90	0.54
2:L:59:PRO:HG2	2:L:62:PHE:CD2	2.43	0.53
3:H:8:GLY:HA3	3:H:20:LEU:CD2	2.37	0.53
2:L:189:HIS:O	2:L:211:PRO:HG2	2.08	0.53
2:L:134:CYS:HB2	2:L:148:TRP:CH2	2.44	0.52
2:L:113:PRO:HB3	2:L:139:PHE:HB3	1.92	0.52
1:A:247:PRO:O	1:A:251:LEU:HG	2.10	0.52
3:H:13:GLN:HA	3:H:112:SER:O	2.10	0.51
1:A:345:GLU:HG3	1:A:432:LEU:HD23	1.91	0.51
3:H:124:LEU:HD11	3:H:143:LEU:HB2	1.92	0.51
1:A:391:TYR:HB3	1:A:410:LEU:HD12	1.92	0.51
2:L:12:SER:HB2	2:L:106(A):LEU:HD21	1.93	0.51
3:H:191:VAL:O	3:H:191:VAL:HG13	2.12	0.50
2:L:190:LYS:O	2:L:211:PRO:HD2	2.11	0.50
2:L:130:ALA:HB3	2:L:181:LEU:HD12	1.92	0.50
1:A:358:MET:O	1:A:414:LYS:HE3	2.11	0.49
1:A:365:LEU:HD21	1:A:417:TRP:CE3	2.47	0.49
2:L:195:GLN:HB3	2:L:206:GLU:HB3	1.94	0.49
2:L:90:VAL:HG12	2:L:97:VAL:HG13	1.92	0.49
2:L:59:PRO:HG2	2:L:62:PHE:HD2	1.78	0.49
1:A:358:MET:SD	1:A:363:VAL:HG11	2.53	0.49
2:L:82:ASP:O	2:L:104:LEU:HD23	2.13	0.48
3:H:94:LYS:HD3	3:H:102:TYR:HD2	1.78	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:22:CYS:HB3	3:H:78:LEU:HB3	1.96	0.47
1:A:365:LEU:HD22	1:A:441:LEU:HD13	1.97	0.47
1:A:425:CYS:O	1:A:438:GLN:HA	2.15	0.47
1:A:367:CYS:HB2	1:A:381:TRP:CZ2	2.50	0.46
2:L:119:PRO:HD2	3:H:125:VAL:O	2.15	0.46
1:A:348:VAL:CG2	1:A:427:VAL:HG11	2.43	0.46
3:H:68:ILE:HB	3:H:81:GLN:HB3	1.97	0.46
1:A:261:CYS:HB3	1:A:304:SER:HB3	1.98	0.46
1:A:370:LYS:HB3	1:A:405:PHE:HB2	1.98	0.46
3:H:142:CYS:HB2	3:H:157:TRP:CZ2	2.51	0.46
2:L:164:PRO:HA	2:L:174:ALA:O	2.16	0.46
1:A:427:VAL:CG1	1:A:437:THR:HG23	2.46	0.46
3:H:20:LEU:HD12	3:H:80:LEU:HD23	1.97	0.45
2:L:195:GLN:CB	2:L:206:GLU:HB3	2.46	0.45
3:H:11:LEU:HD23	3:H:12:VAL:N	2.32	0.45
2:L:14:ALA:HA	2:L:106(A):LEU:HB2	1.97	0.45
2:L:33:VAL:HG21	2:L:71:ALA:HB2	1.98	0.44
2:L:66:ASN:HA	2:L:71:ALA:HA	2.00	0.44
2:L:54:ARG:HD3	2:L:62:PHE:O	2.18	0.44
1:A:422:VAL:HA	1:A:441:LEU:O	2.18	0.44
3:H:67:PHE:CZ	3:H:82:MET:HE3	2.52	0.44
2:L:4:LEU:HA	2:L:24:GLY:O	2.17	0.44
1:A:348:VAL:HG21	1:A:427:VAL:CG1	2.45	0.44
1:A:427:VAL:HG13	1:A:437:THR:CG2	2.48	0.43
2:L:118:PHE:HA	2:L:119:PRO:HD3	1.86	0.43
2:L:33:VAL:HG21	2:L:71:ALA:CB	2.48	0.43
1:A:343:PRO:HB2	1:A:431:ALA:HB2	2.00	0.43
1:A:378:ALA:HB3	1:A:428:MET:HB2	2.01	0.43
2:L:36:TYR:HA	2:L:45:VAL:O	2.19	0.43
1:A:349:TYR:HE2	1:A:370:LYS:HE2	1.84	0.43
2:L:46:LEU:HD11	2:L:49:TYR:HB3	2.01	0.43
1:A:360:LYS:O	1:A:414:LYS:NZ	2.51	0.42
3:H:4:LEU:HD22	3:H:24:THR:HG22	1.99	0.42
2:L:145:THR:OG1	2:L:197:THR:HB	2.20	0.42
2:L:61:ARG:HD2	2:L:76:SER:O	2.19	0.42
3:H:34:MET:HB2	3:H:78:LEU:HD13	2.00	0.42
3:H:2:VAL:HG13	3:H:27:PHE:CD2	2.55	0.42
2:L:13:VAL:HG23	2:L:78:VAL:HG11	2.01	0.42
3:H:188:THR:HG22	3:H:189:SER:N	2.35	0.41
3:H:7:SER:OG	3:H:21:SER:HB2	2.20	0.41
1:A:320:LYS:HA	1:A:334:LYS:O	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:120:PRO:HB3	2:L:131:THR:H	1.85	0.41
3:H:99:VAL:HG23	3:H:100(B):ALA:HB2	2.02	0.41
3:H:12:VAL:O	3:H:111:VAL:HA	2.20	0.41
2:L:62:PHE:HE1	2:L:86:TYR:HE2	1.68	0.41
3:H:48:VAL:HG13	3:H:63:VAL:HG21	2.03	0.41
1:A:346:PRO:HD3	1:A:429:HIS:CD2	2.55	0.40
3:H:2:VAL:HG13	3:H:27:PHE:HD2	1.85	0.40
1:A:377:ILE:HG12	1:A:378:ALA:H	1.87	0.40
1:A:377:ILE:HD11	1:A:427:VAL:HG21	2.04	0.40
1:A:365:LEU:HD21	1:A:417:TRP:CZ3	2.57	0.40
3:H:172:ARG:HB2	3:H:190:GLN:O	2.22	0.40
3:H:6:GLU:HA	3:H:21:SER:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	204/206 (99%)	176 (86%)	23 (11%)	5 (2%)	5	29
2	L	211/213 (99%)	193 (92%)	15 (7%)	3 (1%)	11	43
3	H	223/225 (99%)	185 (83%)	32 (14%)	6 (3%)	5	27
All	All	638/644 (99%)	554 (87%)	70 (11%)	14 (2%)	6	32

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	269	GLU
3	H	150	ASP
1	A	293	GLU
2	L	167	GLN
3	H	180	GLY

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	300	TYR
1	A	339	ALA
2	L	168	SER
3	H	41	PRO
3	H	44	GLY
3	H	130	SER
1	A	297	ASN
3	H	223(A)	PRO
2	L	107	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	194/194 (100%)	184 (95%)	10 (5%)	23	55
2	L	178/178 (100%)	173 (97%)	5 (3%)	43	73
3	H	193/193 (100%)	186 (96%)	7 (4%)	35	67
All	All	565/565 (100%)	543 (96%)	22 (4%)	32	64

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	250	THR
1	A	260	THR
1	A	267	SER
1	A	285	HIS
1	A	294	GLN
1	A	340	LYS
1	A	350	THR
1	A	370	LYS
1	A	414	LYS
1	A	415	SER
2	L	5	THR
2	L	105	THR
2	L	123	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	L	126	GLN
2	L	170	ASN
3	H	39	GLN
3	H	63	VAL
3	H	71	ARG
3	H	107	ILE
3	H	159	TYR
3	H	172	ARG
3	H	178	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	418	GLN
1	A	438	GLN
2	L	171	ASN
2	L	195	GLN
3	H	39	GLN
3	H	82(A)	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	166/206 (80%)	-0.74	0	100 100	2, 21, 67, 90	16 (9%)
2	L	209/213 (98%)	-0.47	0	100 100	2, 54, 54, 72	49 (23%)
3	H	184/225 (81%)	-0.46	3 (1%)	72 59	2, 33, 69, 79	51 (27%)
All	All	559/644 (86%)	-0.55	3 (0%)	91 86	2, 35, 69, 90	116 (20%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	H	118	ALA	2.7
3	H	151	SER	2.4
3	H	223(A)	PRO	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 carbohydrates 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.