



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 09:28 AM EDT

PDB ID : 7WN8 / pdb_00007wn8
Title : Crystal structure of antibody (BC31M5) binds to CD47
Authors : Li, Y.; Wang, W.; Sui, J.; Zhang, S.
Deposited on : 2022-01-17
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	:	4-5-2 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

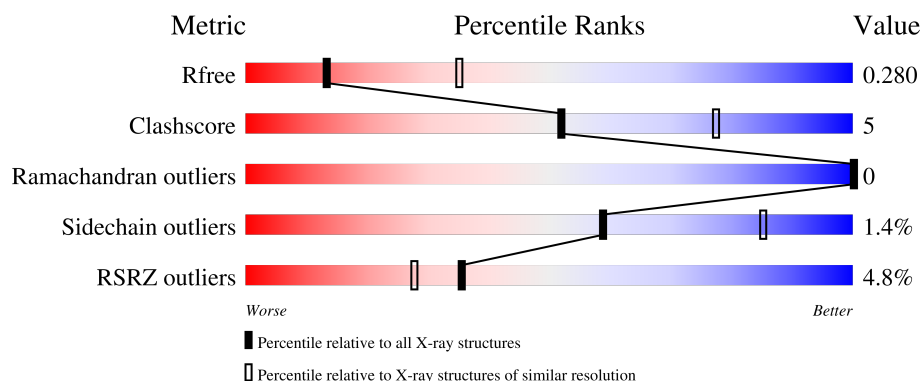
1 Overall quality at a glance

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))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (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 $\leq 5\%$. 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	323	 3% 30% 6% 64%
1	C	323	 6% 27% 5% 68%
2	H	230	 3% 79% 11% 10%
2	P	230	 4% 80% 9% 10%
3	L	215	 90% 9%

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Mol	Chain	Length	Quality of chain
3	Q	215	% 87% 13%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8343 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 Leukocyte surface antigen CD47.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	116	Total	C	N	O	S	0	0	0
			918	583	148	183	4			
1	C	104	Total	C	N	O	S	0	0	0
			814	515	131	164	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	15	GLY	CYS	engineered mutation	UNP Q08722
C	15	GLY	CYS	engineered mutation	UNP Q08722

- Molecule 2 is a protein called BC31M5 Fab Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	207	Total	C	N	O	S	0	0	0
			1553	982	259	305	7			
2	P	206	Total	C	N	O	S	0	0	0
			1551	981	258	305	7			

- Molecule 3 is a protein called BC31M5 Fab Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	214	Total	C	N	O	S	0	0	0
			1658	1032	289	333	4			
3	Q	213	Total	C	N	O	S	0	0	0
			1649	1027	288	330	4			

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	Zn	0	0
			4	4		

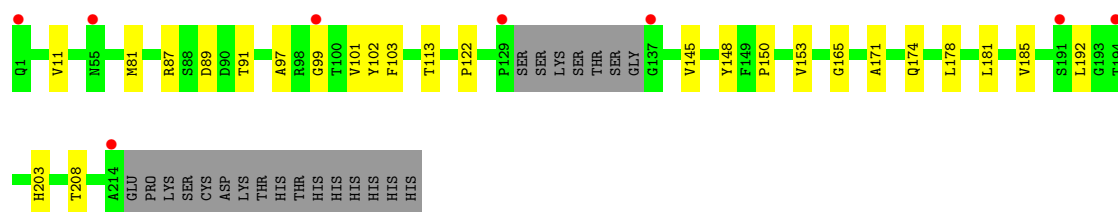
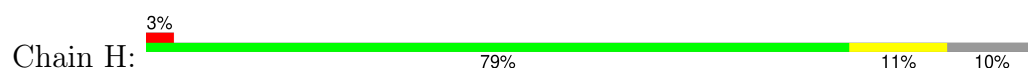
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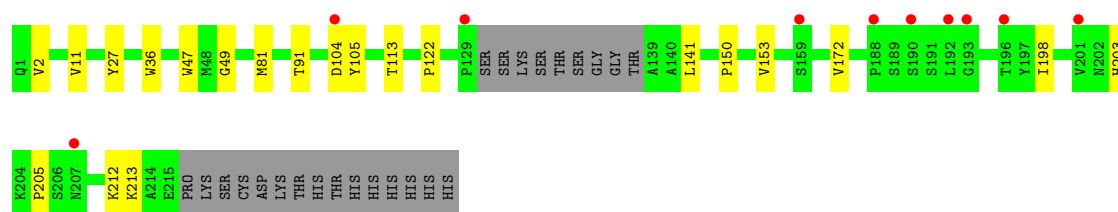
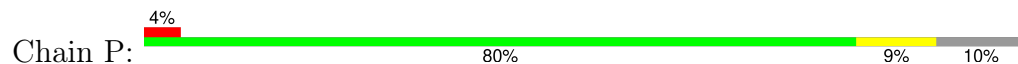
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	2	Total 2	Zn 2	0	0
4	L	1	Total 1	Zn 1	0	0
4	P	1	Total 1	Zn 1	0	0

- Molecule 5 is water.

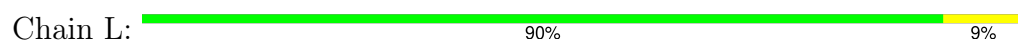
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	23	Total 23	O 23	0	0
5	C	18	Total 18	O 18	0	0
5	H	40	Total 40	O 40	0	0
5	L	36	Total 36	O 36	0	0
5	P	35	Total 35	O 35	0	0
5	Q	40	Total 40	O 40	0	0



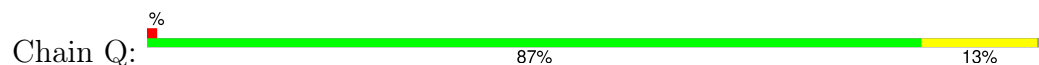
- Molecule 2: BC31M5 Fab Heavy chain



- Molecule 3: BC31M5 Fab Light chain



- Molecule 3: BC31M5 Fab Light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.40Å 94.28Å 174.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.59 – 2.80 49.59 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.59-2.80) 99.8 (49.59-2.80)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.19_4092	Depositor
R, R_{free}	0.243 , 0.276 0.246 , 0.280	Depositor DCC
R_{free} test set	1999 reflections (5.27%)	wwPDB-VP
Wilson B-factor (Å ²)	49.5	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 35.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for k,h,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	8343	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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: ZN

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	0.19	0/933	0.31	0/1261
1	C	0.40	0/825	0.66	1/1114 (0.1%)
2	H	0.18	0/1590	0.35	0/2169
2	P	0.17	0/1588	0.38	0/2166
3	L	0.16	0/1695	0.34	0/2305
3	Q	0.15	0/1686	0.32	0/2293
All	All	0.20	0/8317	0.39	1/11308 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	40	PRO	N-CA-C	8.27	123.29	110.80

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	918	0	909	11	0
1	C	814	0	813	10	0
2	H	1553	0	1520	19	0
2	P	1551	0	1516	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	L	1658	0	1607	15	0
3	Q	1649	0	1601	15	0
4	A	4	0	0	0	0
4	C	2	0	0	0	0
4	L	1	0	0	0	0
4	P	1	0	0	0	0
5	A	23	0	0	0	0
5	C	18	0	0	3	0
5	H	40	0	0	1	0
5	L	36	0	0	1	0
5	P	35	0	0	0	0
5	Q	40	0	0	2	0
All	All	8343	0	7966	82	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 5.

All (82) 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:148:TYR:CE1	2:H:153:VAL:HG23	2.21	0.76
1:C:108:HIS:CE1	5:C:506:HOH:O	2.41	0.72
2:P:198:ILE:HG12	2:P:213:LYS:HG2	1.77	0.65
3:L:13:LEU:HD12	3:L:17:GLU:HB3	1.79	0.64
2:P:104:ASP:OD1	2:P:105:TYR:HD1	1.80	0.64
3:L:17:GLU:HG2	3:Q:24:ARG:HH12	1.63	0.63
3:L:55:ARG:NH2	5:L:401:HOH:O	2.33	0.62
3:L:6:GLN:NE2	3:L:87:TYR:O	2.32	0.61
2:P:104:ASP:OD1	2:P:105:TYR:CD1	2.53	0.61
1:A:55:TYR:HB3	1:A:117:THR:HB	1.82	0.60
1:A:46:MET:HE1	1:A:92:LEU:HA	1.84	0.60
2:H:11:VAL:HG21	2:H:150:PRO:HG3	1.84	0.60
1:C:50:ASN:HA	1:C:92:LEU:HD11	1.82	0.59
2:H:122:PRO:HD2	2:H:208:THR:HG21	1.85	0.58
1:C:108:HIS:HE1	5:C:506:HOH:O	1.83	0.57
2:P:11:VAL:HG21	2:P:150:PRO:HG3	1.87	0.56
3:L:93:THR:O	3:L:96:THR:OG1	2.23	0.56
2:H:97:ALA:HB1	2:H:103:PHE:HB3	1.88	0.55
1:C:40:PRO:HD2	1:C:127:ILE:CD1	2.38	0.53
2:H:148:TYR:CE1	2:H:153:VAL:CG2	2.91	0.53
3:L:61:ASP:N	3:L:61:ASP:OD1	2.37	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:109:ARG:NH1	3:L:110:THR:O	2.42	0.52
2:P:104:ASP:CG	2:P:105:TYR:CD1	2.87	0.52
1:C:73:ASN:ND2	5:C:504:HOH:O	2.43	0.52
3:Q:147:VAL:HG22	3:Q:197:VAL:HG22	1.92	0.51
2:P:212:LYS:HA	2:P:212:LYS:HE2	1.92	0.51
1:A:47:GLU:N	1:A:118:GLU:OE2	2.35	0.51
3:Q:32:ASN:HA	3:Q:51:TRP:HA	1.92	0.51
1:A:37:VAL:HG11	1:A:105:ALA:HB2	1.94	0.50
2:H:122:PRO:HB3	2:H:148:TYR:HB3	1.94	0.49
3:Q:107:ILE:O	3:Q:167:GLN:NE2	2.44	0.49
2:P:91:THR:HG23	2:P:113:THR:HA	1.94	0.49
2:H:203:HIS:HB3	2:H:208:THR:HB	1.94	0.48
3:Q:93:THR:O	3:Q:96:THR:OG1	2.31	0.48
3:Q:95:ASN:OD1	5:Q:301:HOH:O	2.20	0.48
1:C:37:VAL:HG21	1:C:105:ALA:HB2	1.96	0.47
2:H:81:MET:HE3	2:H:81:MET:HB3	1.80	0.47
2:H:174:GLN:HG2	2:H:178:LEU:O	2.15	0.47
3:L:11:LEU:HG	3:L:13:LEU:HD23	1.95	0.47
2:H:171:ALA:HA	2:H:181:LEU:HB3	1.96	0.46
2:H:101:VAL:HG22	3:L:92:HIS:HD2	1.80	0.46
2:H:91:THR:HG23	2:H:113:THR:HA	1.98	0.46
2:H:99:GLY:HA3	2:H:102:TYR:O	2.15	0.46
2:P:212:LYS:HD3	2:P:213:LYS:H	1.81	0.46
2:P:203:HIS:CD2	2:P:205:PRO:HD2	2.51	0.45
1:A:22:PHE:CE1	1:A:43:VAL:HG22	2.52	0.45
2:H:171:ALA:HB2	2:H:181:LEU:HD23	1.98	0.45
3:L:152:ASP:OD2	3:L:190:HIS:ND1	2.29	0.45
1:C:58:TRP:CE2	1:C:98:LEU:HB2	2.52	0.45
1:A:26:LYS:HE2	1:A:26:LYS:HB2	1.73	0.45
3:Q:55:ARG:HD3	3:Q:63:PHE:O	2.17	0.45
2:H:145:VAL:HG11	2:H:153:VAL:HG11	1.99	0.44
2:P:104:ASP:OD1	2:P:105:TYR:N	2.50	0.44
2:H:87:ARG:NH1	5:H:302:HOH:O	2.49	0.44
2:P:36:TRP:CE2	2:P:81:MET:HB2	2.53	0.44
3:Q:143:ARG:NH2	5:Q:302:HOH:O	2.47	0.44
1:A:132:ARG:HG2	1:A:133:VAL:N	2.31	0.44
1:A:46:MET:HG2	1:A:118:GLU:OE2	2.18	0.44
3:Q:19:ALA:HB3	3:Q:76:ILE:HB	2.00	0.44
3:L:18:ARG:HE	3:L:18:ARG:HB2	1.62	0.44
2:P:2:VAL:HG13	2:P:27:TYR:CD1	2.53	0.44
2:P:122:PRO:HD3	2:P:203:HIS:ND1	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:146:LYS:HB3	3:Q:198:THR:HB	1.99	0.44
1:C:63:ARG:HH12	1:C:80:ASP:CG	2.26	0.43
1:A:43:VAL:HG21	1:A:116:VAL:HG11	2.00	0.43
1:A:58:TRP:CH2	1:A:114:CYS:HB2	2.53	0.43
2:H:87:ARG:HD2	2:H:89:ASP:HB2	2.01	0.43
3:L:106:GLU:HG2	3:L:167:GLN:OE1	2.19	0.42
3:L:109:ARG:HG3	3:L:110:THR:O	2.20	0.41
3:L:145:ALA:HB2	3:L:199:HIS:HD2	1.84	0.41
3:Q:38:GLN:HB2	3:Q:48:LEU:HD11	2.01	0.41
1:C:58:TRP:HE1	1:C:86:ILE:HD11	1.85	0.41
3:Q:141:TYR:CG	3:Q:142:PRO:HA	2.55	0.41
2:H:165:GLY:O	2:H:185:VAL:HA	2.20	0.41
3:L:121:PRO:HB2	3:L:126:LEU:HD21	2.02	0.41
3:Q:109:ARG:HD2	3:Q:171:ASP:O	2.21	0.41
1:A:58:TRP:CE2	1:A:98:LEU:HB2	2.56	0.41
3:Q:21:LEU:HD23	3:Q:103:THR:HB	2.03	0.41
2:H:192:LEU:HD23	2:H:192:LEU:HA	1.87	0.40
2:P:47:TRP:CH2	2:P:49:GLY:HA2	2.57	0.40
3:Q:141:TYR:CD1	3:Q:142:PRO:HA	2.56	0.40
1:C:66:TYR:CE1	1:C:86:ILE:HG13	2.57	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	114/323 (35%)	112 (98%)	2 (2%)	0	100	100
1	C	100/323 (31%)	98 (98%)	2 (2%)	0	100	100
2	H	203/230 (88%)	200 (98%)	3 (2%)	0	100	100
2	P	202/230 (88%)	200 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	L	212/215 (99%)	206 (97%)	6 (3%)	0	100	100
3	Q	211/215 (98%)	205 (97%)	6 (3%)	0	100	100
All	All	1042/1536 (68%)	1021 (98%)	21 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	105/274 (38%)	101 (96%)	4 (4%)	29	64
1	C	94/274 (34%)	91 (97%)	3 (3%)	34	70
2	H	172/194 (89%)	172 (100%)	0	100	100
2	P	172/194 (89%)	169 (98%)	3 (2%)	53	83
3	L	188/189 (100%)	187 (100%)	1 (0%)	81	93
3	Q	187/189 (99%)	185 (99%)	2 (1%)	65	87
All	All	918/1314 (70%)	905 (99%)	13 (1%)	59	85

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	VAL
1	A	38	VAL
1	A	107	SER
1	A	134	VAL
1	C	36	THR
1	C	77	VAL
1	C	103	SER
3	L	106	GLU
2	P	141	LEU
2	P	153	VAL
2	P	172	VAL
3	Q	34	LEU

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Mol	Chain	Res	Type
3	Q	98	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
3	L	91	GLN
2	P	167	HIS
3	Q	161	GLN

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](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

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	116/323 (35%)	0.81	11 (9%) 14 10	33, 53, 71, 74	0
1	C	104/323 (32%)	1.11	19 (18%) 3 2	37, 60, 79, 81	0
2	H	207/230 (90%)	0.26	8 (3%) 43 34	28, 36, 57, 69	0
2	P	206/230 (89%)	0.53	10 (4%) 35 27	31, 45, 64, 76	0
3	L	214/215 (99%)	0.16	1 (0%) 87 82	28, 38, 46, 52	0
3	Q	213/215 (99%)	0.20	2 (0%) 81 74	31, 41, 49, 59	0
All	All	1060/1536 (69%)	0.42	51 (4%) 35 28	28, 41, 68, 81	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	132	ARG	4.9
1	C	108	HIS	4.4
1	C	107	SER	4.0
1	A	134	VAL	3.9
1	A	32	PHE	3.8
2	H	191	SER	3.7
1	A	133	VAL	3.5
1	C	109	THR	3.5
1	C	110	GLY	3.5
2	P	190	SER	3.2
2	H	55	ASN	3.0
2	H	194	THR	3.0
2	P	104	ASP	2.9
1	C	28	VAL	2.8
1	A	49	GLN	2.7
3	Q	1	GLU	2.7
2	P	129	PRO	2.6
1	A	31	THR	2.6
1	A	35	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
2	P	201	VAL	2.6
2	H	1	GLN	2.6
3	L	77	SER	2.5
1	A	79	THR	2.5
1	A	63	ARG	2.5
1	C	129	LEU	2.4
2	P	192	LEU	2.4
1	C	78	PRO	2.4
1	C	79	THR	2.4
1	C	127	ILE	2.4
2	H	129	PRO	2.4
1	A	33	CYS	2.4
1	A	78	PRO	2.4
2	H	214	ALA	2.3
1	C	103	SER	2.3
2	H	137	GLY	2.3
1	C	36	THR	2.3
2	P	196	THR	2.3
1	C	37	VAL	2.2
3	Q	28	SER	2.2
2	P	159	SER	2.2
1	C	102	LYS	2.2
1	C	60	PHE	2.2
1	C	49	GLN	2.1
2	P	207	ASN	2.1
2	P	188	PRO	2.1
2	P	193	GLY	2.1
2	H	99	GLY	2.1
1	C	76	THR	2.1
1	C	62	GLY	2.0
1	C	104	ASP	2.0
1	C	25	THR	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	ZN	A	402	1/1	0.61	0.30	214,214,214,214	0
4	ZN	P	301	1/1	0.61	0.22	58,58,58,58	0
4	ZN	A	403	1/1	0.75	0.23	168,168,168,168	0
4	ZN	L	301	1/1	0.94	0.12	56,56,56,56	0
4	ZN	C	402	1/1	0.95	0.07	34,34,34,34	0
4	ZN	C	401	1/1	0.96	0.07	43,43,43,43	0
4	ZN	A	404	1/1	0.98	0.04	30,30,30,30	0
4	ZN	A	401	1/1	1.00	0.07	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.