



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 18, 2022 – 02:05 PM EDT

PDB ID : 7RDV
Title : TFH TCR bound to MHC Class II IAd presenting aggrecan epitope
Authors : Lim, J.J.; Rossjohn, J.; Reid, H.
Deposited on : 2021-07-11
Resolution : 2.90 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	:	2.29
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)	:	2.29

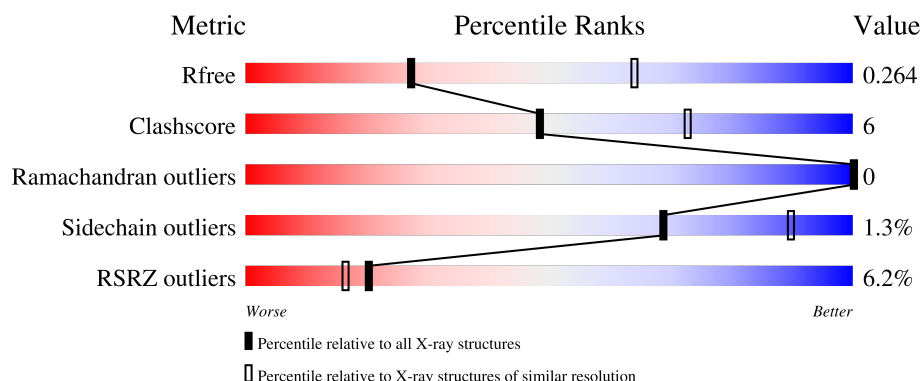
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	181	<div> <div>9%</div> <div>82%</div> <div>18%</div> </div>
2	B	186	<div> <div>12%</div> <div>75%</div> <div>17%</div> <div>8%</div> </div>
3	C	205	<div> <div>2%</div> <div>75%</div> <div>14%</div> <div>11%</div> </div>
4	D	240	<div> <div>2%</div> <div>88%</div> <div>12%</div> </div>
5	H	12	<div> <div>75%</div> <div>17%</div> <div>8%</div> </div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6052 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 H-2 class II histocompatibility antigen, A-D alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	181	Total	C	N	O	S	0	0	0
			1393	900	225	266	2			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	179	THR	-	expression tag	UNP P04228
A	180	SER	-	expression tag	UNP P04228
A	181	GLY	-	expression tag	UNP P04228

- Molecule 2 is a protein called H-2 class II histocompatibility antigen, A-D beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	171	Total	C	N	O	S	0	0	0
			1350	844	239	262	5			

- Molecule 3 is a protein called TFH TCR alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	183	Total	C	N	O	S	0	0	0
			1371	855	229	279	8			

- Molecule 4 is a protein called TFH TCR beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	240	Total	C	N	O	S	0	0	0
			1808	1138	324	340	6			

- Molecule 5 is a protein called AggreCAN core peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	H	12	Total	C	N	O	0	0	0
			95	56	20	19			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			14	8	1	5		
6	A	1	Total	C	N	O	0	0
			14	8	1	5		

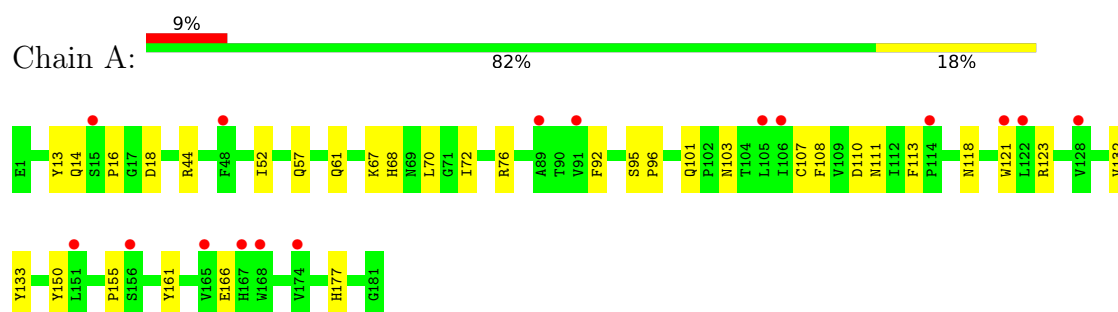
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total O 1 1	0	0
7	B	2	Total O 2 2	0	0
7	D	4	Total O 4 4	0	0

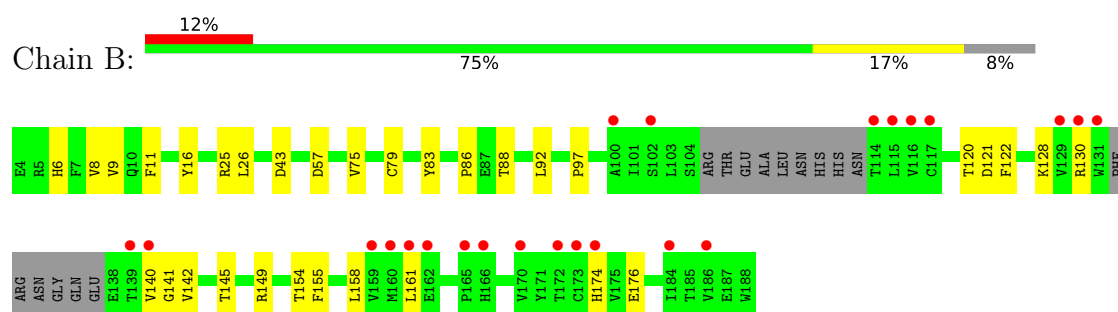
3 Residue-property plots [i](#)

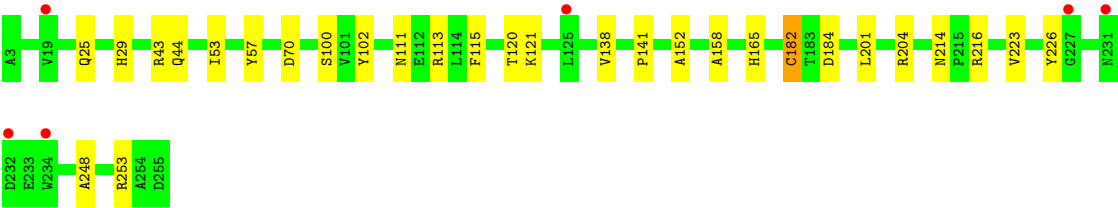
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: H-2 class II histocompatibility antigen, A-D alpha chain



- Molecule 2: H-2 class II histocompatibility antigen, A-D beta chain





● Molecule 5: Aggrecan core peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	102.56Å 102.56Å 195.69Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.61 – 2.90 49.61 – 2.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.61-2.90) 100.0 (49.61-2.90)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.54 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.237 , 0.265 0.237 , 0.264	Depositor DCC
R_{free} test set	1359 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	84.7	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 59.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6052	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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](#)

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

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	0.24	0/1437	0.43	0/1972
2	B	0.24	0/1379	0.44	0/1878
3	C	0.28	0/1396	0.47	0/1901
4	D	0.26	0/1859	0.43	0/2540
5	H	0.25	0/95	0.39	0/126
All	All	0.26	0/6166	0.44	0/8417

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 [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	1393	0	1271	22	0
2	B	1350	0	1238	23	0
3	C	1371	0	1233	19	0
4	D	1808	0	1645	18	0
5	H	95	0	90	2	0
6	A	28	0	26	0	0
7	A	1	0	0	0	0
7	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	D	4	0	0	0	0
All	All	6052	0	5503	69	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 6.

All (69) 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:C:31:ASN:HA	3:C:58:SER:HB3	1.75	0.68
3:C:57:ARG:NH1	4:D:111:ASN:O	2.29	0.65
1:A:16:PRO:HG2	2:B:6:HIS:HA	1.77	0.65
3:C:40:PRO:HG2	3:C:105:ALA:HB3	1.79	0.65
1:A:118:ASN:HB2	1:A:166:GLU:HB2	1.81	0.62
3:C:32:TYR:HB3	3:C:107:SER:HB3	1.83	0.61
1:A:14:GLN:HG2	2:B:8:VAL:HG22	1.84	0.59
1:A:70:LEU:HD13	2:B:9:VAL:HG13	1.86	0.58
1:A:61:GLN:OE1	5:H:93:ARG:NH2	2.24	0.57
4:D:214:ASN:HD21	4:D:216:ARG:NH1	2.01	0.57
2:B:149:ARG:HG3	2:B:155:PHE:HE1	1.70	0.57
3:C:107:SER:HB2	3:C:114:ILE:HG22	1.86	0.57
2:B:130:ARG:HG3	2:B:174:HIS:HB3	1.88	0.55
2:B:142:VAL:HG22	2:B:161:LEU:HD13	1.88	0.55
2:B:83:TYR:O	2:B:88:THR:OG1	2.19	0.54
2:B:140:VAL:HG12	2:B:141:GLY:H	1.71	0.54
4:D:25:GLN:NE2	4:D:29:HIS:O	2.40	0.53
3:C:150:LEU:HG	3:C:152:THR:HG23	1.90	0.53
4:D:165:HIS:HB3	4:D:226:TYR:HB2	1.91	0.53
1:A:110:ASP:OD1	1:A:111:ASN:N	2.41	0.52
2:B:97:PRO:HB3	2:B:122:PHE:HB3	1.92	0.52
3:C:6:GLN:HG2	3:C:120:THR:HG23	1.91	0.52
3:C:44:GLN:OE1	4:D:44:GLN:NE2	2.33	0.51
4:D:141:PRO:HG2	4:D:152:ALA:HB1	1.92	0.51
2:B:128:LYS:HB3	2:B:176:GLU:HB3	1.93	0.51
2:B:121:ASP:HA	2:B:154:THR:HB	1.91	0.51
1:A:57:GLN:HG2	3:C:112:ASN:ND2	2.27	0.50
1:A:13:TYR:CZ	1:A:67:LYS:HG3	2.46	0.50
1:A:95:SER:O	1:A:103:ASN:ND2	2.39	0.50
2:B:16:TYR:HB2	2:B:25:ARG:HB3	1.92	0.50
1:A:61:GLN:HB2	3:C:112:ASN:HD21	1.78	0.49
3:C:53:ILE:HG13	3:C:54:ILE:HG22	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:43:ARG:HB3	4:D:53:ILE:HD11	1.96	0.48
3:C:176:LEU:HB3	4:D:182:CYS:HB2	1.95	0.48
1:A:44:ARG:HH22	1:A:133:TYR:HE1	1.62	0.48
1:A:13:TYR:OH	1:A:18:ASP:OD1	2.16	0.48
1:A:123:ARG:HA	1:A:161:TYR:HD1	1.79	0.48
1:A:61:GLN:HG3	4:D:57:TYR:HB2	1.96	0.48
1:A:68:HIS:O	1:A:72:ILE:HG12	2.14	0.47
4:D:184:ASP:OD1	4:D:204:ARG:NH2	2.49	0.46
1:A:52:ILE:HD13	2:B:86:PRO:HB3	1.96	0.46
4:D:102:TYR:HB2	4:D:120:THR:HG23	1.97	0.46
4:D:138:VAL:HG23	4:D:248:ALA:HB3	1.98	0.45
2:B:25:ARG:NE	2:B:43:ASP:OD2	2.50	0.45
1:A:92:PHE:HE1	1:A:108:PHE:HB2	1.81	0.45
2:B:26:LEU:HD23	2:B:79:CYS:SG	2.56	0.44
4:D:158:ALA:HB2	4:D:223:VAL:HG21	1.99	0.44
4:D:43:ARG:HH21	4:D:100:SER:HB2	1.83	0.44
2:B:11:PHE:CD2	5:H:96:VAL:HG22	2.53	0.44
2:B:128:LYS:HE3	2:B:128:LYS:HB2	1.84	0.43
1:A:107:CYS:HB2	1:A:121:TRP:CZ2	2.53	0.43
3:C:6:GLN:NE2	3:C:120:THR:HG23	2.34	0.43
3:C:199:CYS:HA	3:C:202:ALA:HB2	1.98	0.43
4:D:113:ARG:HD2	4:D:115:PHE:CZ	2.54	0.43
1:A:101:GLN:O	1:A:155:PRO:HD2	2.19	0.42
1:A:96:PRO:HD3	2:B:120:THR:HG21	2.01	0.42
1:A:132:VAL:HA	1:A:150:TYR:O	2.19	0.42
2:B:140:VAL:HG12	2:B:141:GLY:N	2.35	0.42
2:B:145:THR:HG22	2:B:158:LEU:H	1.84	0.42
3:C:6:GLN:HG3	3:C:23:CYS:SG	2.60	0.42
3:C:113:ARG:HH12	4:D:70:ASP:CG	2.24	0.41
3:C:163:SER:HB3	3:C:170:ILE:HG13	2.02	0.41
3:C:7:LEU:O	3:C:9:SER:N	2.54	0.41
4:D:121:LYS:HG2	4:D:165:HIS:CE1	2.55	0.41
1:A:76:ARG:HH12	2:B:57:ASP:CG	2.24	0.41
2:B:75:VAL:O	2:B:79:CYS:HB2	2.21	0.41
2:B:83:TYR:CZ	2:B:92:LEU:HD21	2.56	0.41
3:C:79:LEU:HD22	3:C:89:LEU:HD11	2.02	0.41
4:D:184:ASP:HB2	4:D:201:LEU:HD12	2.02	0.41

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	179/181 (99%)	172 (96%)	7 (4%)	0	100	100
2	B	165/186 (89%)	156 (94%)	9 (6%)	0	100	100
3	C	177/205 (86%)	173 (98%)	4 (2%)	0	100	100
4	D	238/240 (99%)	227 (95%)	11 (5%)	0	100	100
5	H	10/12 (83%)	10 (100%)	0	0	100	100
All	All	769/824 (93%)	738 (96%)	31 (4%)	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	146/162 (90%)	144 (99%)	2 (1%)	67	89
2	B	141/170 (83%)	141 (100%)	0	100	100
3	C	145/187 (78%)	143 (99%)	2 (1%)	67	89
4	D	182/206 (88%)	180 (99%)	2 (1%)	73	92
5	H	10/10 (100%)	8 (80%)	2 (20%)	1	4
All	All	624/735 (85%)	616 (99%)	8 (1%)	69	90

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	PHE
1	A	177	HIS
3	C	104	CYS
3	C	111	ASN
4	D	182	CYS
4	D	253	ARG
5	H	93	ARG
5	H	95	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
3	C	112	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	NAG	A	201	1	14,14,15	0.27	0	17,19,21	0.47	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	A	202	1	14,14,15	0.23	0	17,19,21	0.48	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	A	201	1	-	0/6/23/26	0/1/1/1
6	NAG	A	202	1	-	0/6/23/26	0/1/1/1

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 [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	181/181 (100%)	0.63	16 (8%)	10	7	76, 109, 161, 189	0
2	B	171/186 (91%)	0.74	23 (13%)	3	2	68, 106, 165, 194	0
3	C	183/205 (89%)	0.23	4 (2%)	62	59	58, 83, 135, 147	0
4	D	240/240 (100%)	0.07	6 (2%)	57	55	64, 90, 142, 176	0
5	H	12/12 (100%)	0.32	0	100	100	67, 76, 96, 106	0
All	All	787/824 (95%)	0.39	49 (6%)	20	16	58, 97, 156, 194	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	116	VAL	7.3
2	B	165	PRO	4.9
4	D	231	ASN	4.4
2	B	102	SER	4.3
2	B	129	VAL	4.2
2	B	174	HIS	3.9
2	B	139	THR	3.9
4	D	125	LEU	3.8
2	B	172	THR	3.7
1	A	151	LEU	3.6
2	B	131	TRP	3.6
2	B	100	ALA	3.5
2	B	130	ARG	3.4
1	A	91	VAL	3.2
1	A	174	VAL	3.1
1	A	121	TRP	3.1
2	B	161	LEU	3.1
2	B	159	VAL	3.0
2	B	166	HIS	2.9
1	A	105	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	115	LEU	2.8
2	B	117	CYS	2.8
1	A	89	ALA	2.7
1	A	168	TRP	2.7
2	B	186	VAL	2.7
3	C	205	ASN	2.6
1	A	165	VAL	2.6
4	D	232	ASP	2.5
3	C	200	ALA	2.5
4	D	234	TRP	2.5
3	C	176	LEU	2.5
2	B	162	GLU	2.4
3	C	7	LEU	2.4
2	B	170	VAL	2.3
2	B	160	MET	2.3
1	A	114	PRO	2.2
1	A	48	PHE	2.2
1	A	15	SER	2.2
1	A	106	ILE	2.2
2	B	140	VAL	2.1
1	A	128	VAL	2.1
1	A	156	SER	2.1
2	B	173	CYS	2.1
1	A	122	LEU	2.1
1	A	167	HIS	2.1
2	B	184	ILE	2.1
4	D	19	VAL	2.0
4	D	227	GLY	2.0
2	B	114	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 monosaccharides 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(\AA^2)	Q<0.9
6	NAG	A	201	14/15	0.74	0.36	139,139,139,139	0
6	NAG	A	202	14/15	0.74	0.20	121,121,121,121	0

6.5 Other polymers [i](#)

There are no such residues in this entry.