



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:27 AM UTC

PDB ID : 3QDG / pdb_00003qdg
Title : The complex between TCR DMF5 and human Class I MHC HLA-A2 with the bound MART-1(26-35)(A27L) peptide
Authors : Borbulevych, O.Y.; Baker, B.M.
Deposited on : 2011-01-18
Resolution : 2.69 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

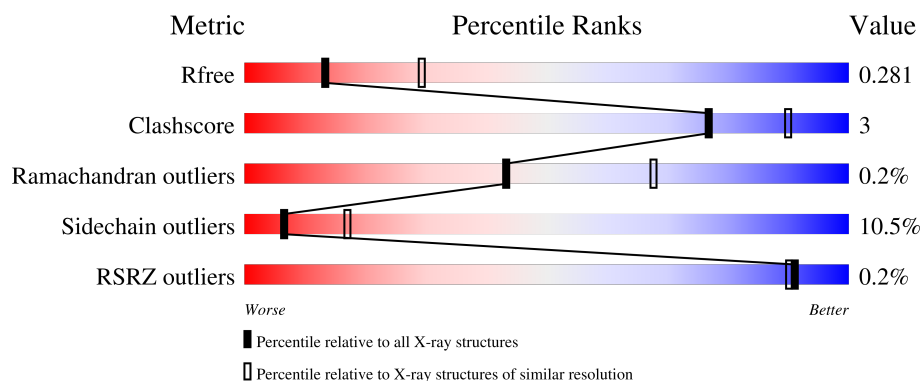
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	275	
2	B	100	
3	C	10	
4	D	199	
5	E	242	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6633 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 HLA class I histocompatibility antigen, A-2 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	1	0
			2255	1408	412	426	9			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	1	0
			842	537	141	159	5			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769

- Molecule 3 is a protein called Mart-1 (26-35) peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	10	Total	C	N	O	0	0	0
			69	45	10	14			

- Molecule 4 is a protein called DMF5 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	199	Total	C	N	O	S	0	0	0
			1546	965	255	318	8			

- Molecule 5 is a protein called DMF5 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	242	Total	C	N	O	S	0	0	0
			1897	1191	333	365	8			

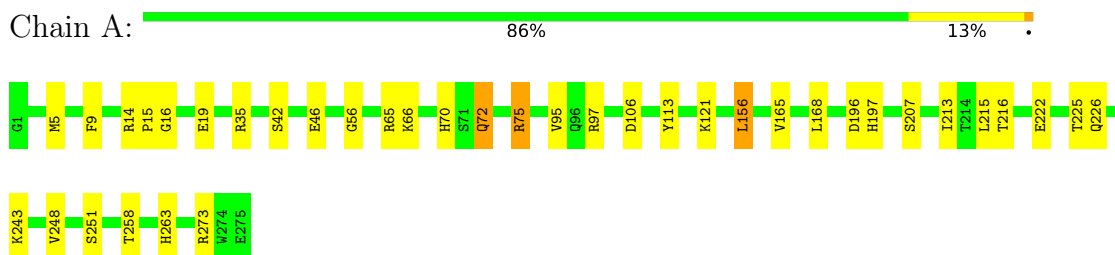
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	9	Total 9	O 9	0	0
6	B	8	Total 8	O 8	0	0
6	C	3	Total 3	O 3	0	0
6	D	1	Total 1	O 1	0	0
6	E	3	Total 3	O 3	0	0

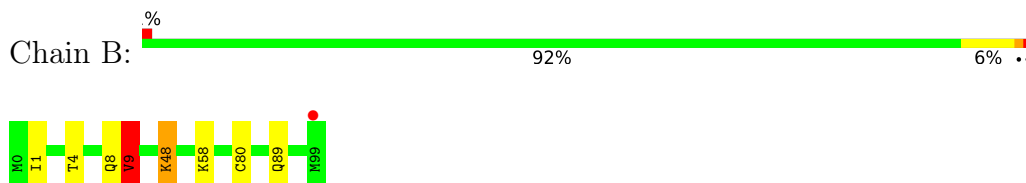
3 Residue-property plots

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.

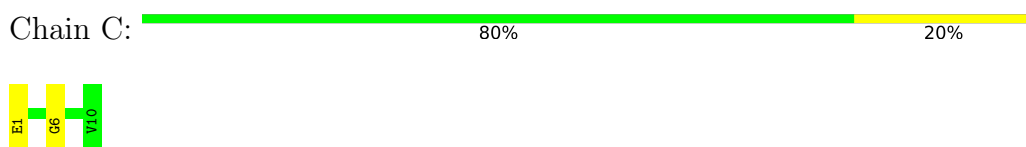
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



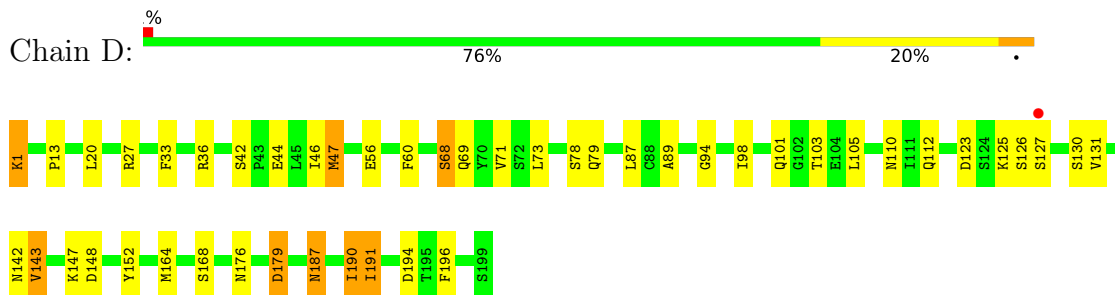
- Molecule 2: Beta-2-microglobulin



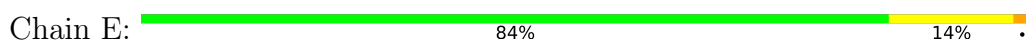
- Molecule 3: Mart-1 (26-35) peptide

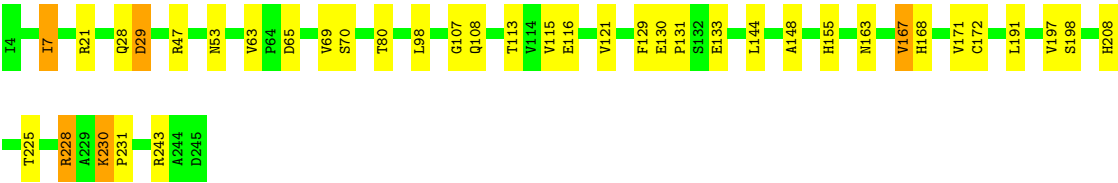


- Molecule 4: DMF5 alpha chain



- Molecule 5: DMF5 beta chain





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	228.36Å 46.55Å 86.04Å 90.00° 106.73° 90.00°	Depositor
Resolution (Å)	20.00 – 2.69 20.00 – 2.69	Depositor EDS
% Data completeness (in resolution range)	90.2 (20.00-2.69) 89.9 (20.00-2.69)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.80 (at 2.71Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.218 , 0.278 0.217 , 0.281	Depositor DCC
R_{free} test set	1123 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	66.4	Xtriage
Anisotropy	0.236	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 66.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6633	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.94% 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](#)

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	0.58	0/2323	1.04	9/3151 (0.3%)
2	B	0.58	0/868	0.96	1/1169 (0.1%)
3	C	0.72	0/68	0.98	0/90
4	D	0.58	0/1579	0.97	5/2135 (0.2%)
5	E	0.52	0/1946	0.93	3/2651 (0.1%)
All	All	0.57	0/6784	0.98	18/9196 (0.2%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	16	GLY	N-CA-C	-9.54	101.98	115.43
1	A	56	GLY	CA-C-N	6.45	126.38	119.28
1	A	56	GLY	C-N-CA	6.45	126.38	119.28
1	A	14	ARG	CA-C-N	-6.34	111.92	119.84
1	A	14	ARG	C-N-CA	-6.34	111.92	119.84
4	D	176	ASN	N-CA-C	5.85	116.66	108.00
4	D	78	SER	N-CA-C	5.81	118.67	109.96
5	E	230	LYS	CA-C-N	5.71	126.98	119.84
5	E	230	LYS	C-N-CA	5.71	126.98	119.84
1	A	46	GLU	CA-C-N	5.57	125.50	119.76
1	A	46	GLU	C-N-CA	5.57	125.50	119.76
1	A	42	SER	N-CA-C	5.56	118.27	111.82
4	D	147	LYS	N-CA-C	-5.55	105.61	112.72
1	A	251	SER	N-CA-C	5.24	117.83	110.23
2	B	9	VAL	CB-CA-C	-5.10	103.23	110.83
5	E	167	VAL	N-CA-C	5.09	115.81	108.48
4	D	143	VAL	N-CA-C	5.06	115.41	108.12
4	D	190	ILE	N-CA-C	5.01	116.85	109.63

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	2255	0	2109	11	0
2	B	842	0	812	2	0
3	C	69	0	79	1	0
4	D	1546	0	1460	14	1
5	E	1897	0	1808	15	0
6	A	9	0	0	0	0
6	B	8	0	0	0	0
6	C	3	0	0	0	0
6	D	1	0	0	0	0
6	E	3	0	0	0	0
All	All	6633	0	6268	37	1

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 3.

All (37) 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)
5:E:130:GLU:OE2	5:E:243:ARG:NH1	2.21	0.74
5:E:121:VAL:O	5:E:228:ARG:NH2	2.26	0.68
1:A:19:GLU:OE2	1:A:75:ARG:NH2	2.32	0.62
1:A:72:GLN:OE1	5:E:53:ASN:ND2	2.30	0.61
1:A:66:LYS:O	1:A:70:HIS:ND1	2.36	0.57
4:D:42:SER:OG	5:E:107:GLY:O	2.22	0.57
4:D:33:PHE:HB2	4:D:89:ALA:HB3	1.90	0.53
5:E:230:LYS:HA	5:E:231:PRO:HD3	1.68	0.47
1:A:65:ARG:HB3	4:D:94:GLY:H	1.78	0.47
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.96	0.47
1:A:258:THR:HG22	1:A:273:ARG:HB3	1.96	0.46
5:E:131:PRO:HD3	5:E:144:LEU:HG	1.97	0.46
4:D:142:ASN:OD1	4:D:142:ASN:N	2.48	0.46
4:D:179:ASP:OD1	4:D:179:ASP:N	2.44	0.46
2:B:48:LYS:H	2:B:48:LYS:HG3	1.30	0.46
5:E:113:THR:OG1	5:E:155:HIS:NE2	2.39	0.45
4:D:187:ASN:HD22	4:D:191:ILE:HG12	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:LEU:HD22	1:A:243:LYS:HD3	1.99	0.44
4:D:60:PHE:HD1	4:D:73:LEU:HD11	1.82	0.44
4:D:36:ARG:NH2	4:D:44:GLU:OE2	2.50	0.44
4:D:1:LYS:HE3	4:D:1:LYS:HB2	1.75	0.44
4:D:194:ASP:OD2	4:D:194:ASP:N	2.50	0.44
1:A:156:LEU:HD13	3:C:6:GLY:HA3	1.99	0.44
4:D:46:ILE:HG22	4:D:47:MET:HE2	2.00	0.44
5:E:108:GLN:H	5:E:108:GLN:HG2	1.59	0.43
4:D:164:MET:HE1	5:E:198:SER:HB3	2.00	0.43
2:B:9:VAL:HG13	2:B:80:CYS:HB2	2.01	0.43
4:D:191:ILE:H	4:D:191:ILE:HG13	1.58	0.42
5:E:7:ILE:HG23	5:E:7:ILE:HD13	1.68	0.42
1:A:213:ILE:HG13	1:A:263:HIS:HB2	2.01	0.42
5:E:29:ASP:OD2	5:E:29:ASP:N	2.53	0.41
1:A:9:PHE:HB2	1:A:97:ARG:HB3	2.03	0.41
5:E:7:ILE:HD12	5:E:7:ILE:HG21	1.68	0.41
5:E:148:ALA:HB3	5:E:191:LEU:HB3	2.02	0.41
4:D:123:ASP:HA	5:E:129:PHE:HA	2.03	0.41
1:A:72:GLN:H	1:A:72:GLN:HG2	1.71	0.40
5:E:163:ASN:HA	5:E:208:HIS:HB3	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:68:SER:OG	4:D:110:ASN:OD1[2_657]	2.17	0.03

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	274/275 (100%)	265 (97%)	7 (3%)	2 (1%)	18	41

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
3	C	8/10 (80%)	8 (100%)	0	0	100	100
4	D	197/199 (99%)	179 (91%)	18 (9%)	0	100	100
5	E	240/242 (99%)	227 (95%)	13 (5%)	0	100	100
All	All	817/826 (99%)	775 (95%)	40 (5%)	2 (0%)	43	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	15	PRO
1	A	226	GLN

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	232/231 (100%)	216 (93%)	16 (7%)	14	34
2	B	96/95 (101%)	89 (93%)	7 (7%)	13	32
3	C	7/7 (100%)	6 (86%)	1 (14%)	3	8
4	D	176/176 (100%)	146 (83%)	30 (17%)	2	6
5	E	204/204 (100%)	183 (90%)	21 (10%)	7	18
All	All	715/713 (100%)	640 (90%)	75 (10%)	6	17

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	ARG
1	A	72	GLN
1	A	75	ARG
1	A	95	VAL
1	A	106	ASP
1	A	113	TYR

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Mol	Chain	Res	Type
1	A	121	LYS
1	A	156	LEU
1	A	165	VAL
1	A	196	ASP
1	A	197	HIS
1	A	207	SER
1	A	216	THR
1	A	222	GLU
1	A	225	THR
1	A	248	VAL
2	B	1	ILE
2	B	4	THR
2	B	8	GLN
2	B	9	VAL
2	B	48	LYS
2	B	58	LYS
2	B	89	GLN
3	C	1	GLU
4	D	1	LYS
4	D	13	PRO
4	D	20	LEU
4	D	27	ARG
4	D	47	MET
4	D	56	GLU
4	D	68	SER
4	D	69	GLN
4	D	71	VAL
4	D	79	GLN
4	D	87	LEU
4	D	98	ILE
4	D	101	GLN
4	D	103	THR
4	D	105	LEU
4	D	112	GLN
4	D	125	LYS
4	D	126	SER
4	D	127	SER
4	D	130	SER
4	D	131	VAL
4	D	143	VAL
4	D	148	ASP
4	D	152	TYR

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Mol	Chain	Res	Type
4	D	168	SER
4	D	179	ASP
4	D	187	ASN
4	D	190	ILE
4	D	191	ILE
4	D	196	PHE
5	E	7	ILE
5	E	21	ARG
5	E	28	GLN
5	E	29	ASP
5	E	47	ARG
5	E	63	VAL
5	E	65	ASP
5	E	69	VAL
5	E	70	SER
5	E	80	THR
5	E	98	LEU
5	E	115	VAL
5	E	116	GLU
5	E	133	GLU
5	E	167	VAL
5	E	168	HIS
5	E	171	VAL
5	E	172	CYS
5	E	197	VAL
5	E	225	THR
5	E	228	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
2	B	13	HIS
4	D	65	ASN
4	D	69	GLN
4	D	187	ASN
4	D	188	ASN
5	E	53	ASN
5	E	108	GLN
5	E	176	GLN
5	E	208	HIS
5	E	226	GLN

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Mol	Chain	Res	Type
5	E	234	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](#)

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 [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	275/275 (100%)	-0.42	0 100 100	42, 74, 122, 146	1 (0%)
2	B	100/100 (100%)	-0.53	1 (1%) 79 78	52, 66, 87, 97	1 (1%)
3	C	10/10 (100%)	-0.61	0 100 100	55, 62, 66, 75	0
4	D	199/199 (100%)	-0.00	1 (0%) 87 86	66, 106, 159, 170	0
5	E	242/242 (100%)	-0.17	0 100 100	61, 102, 137, 151	0
All	All	826/826 (100%)	-0.26	2 (0%) 91 90	42, 86, 145, 170	2 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	127	SER	3.5
2	B	99[A]	MET	3.1

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.