



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:33 AM UTC

PDB ID : 6G3J / pdb_00006g3j
Title : MHC A02 Allele presenting MTSAIGILVP
Authors : Rizkallah, P.J.; Sewell, A.K.
Deposited on : 2018-03-26
Resolution : 2.45 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

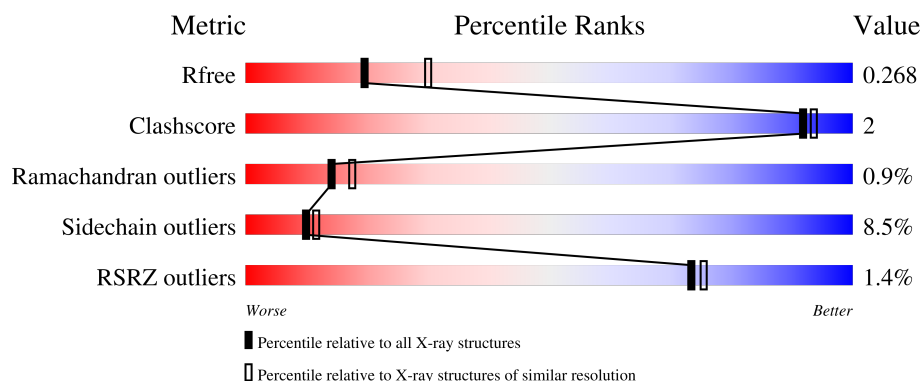
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	276	86% 13%
1	D	276	86% 13%
2	B	100	83% 16%
2	E	100	85% 14%
3	C	10	90% 10%

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Mol	Chain	Length	Quality of chain
3	F	10	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: a green segment representing 70%, a yellow segment representing 20%, and an orange segment representing 10%. The percentages are labeled below the corresponding segments.

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6354 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	276	Total	C	N	O	S	0	1	0
			2264	1414	413	428	9			
1	D	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	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	0	0
			837	533	141	159	4			
2	E	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			

There are 2 discrepancies between the modelled and reference sequences:

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

- Molecule 3 is a protein called MET-THR-SER-ALA-ILE-GLY-ILE-LEU-PRO-VAL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	10	Total	C	N	O	S	0	0	0
			69	45	10	13	1			
3	F	10	Total	C	N	O	S	0	0	0
			69	45	10	13	1			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	O	0	0
			6	6		

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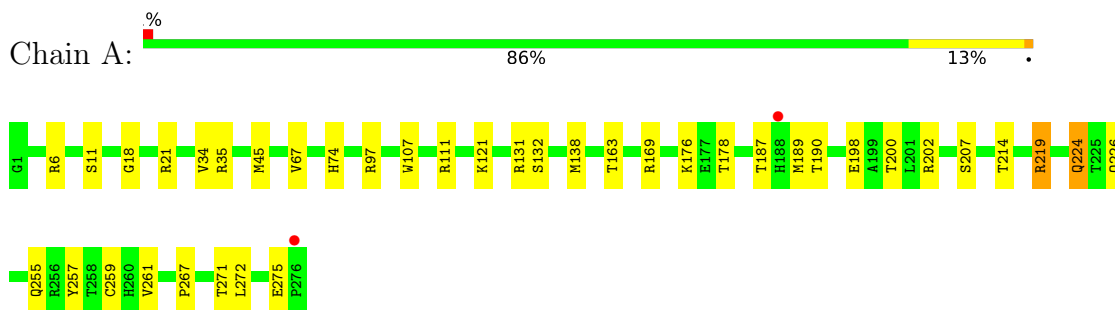
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	4	Total 4	O 4	0	0
4	C	1	Total 1	O 1	0	0
4	D	6	Total 6	O 6	0	0
4	E	7	Total 7	O 7	0	0

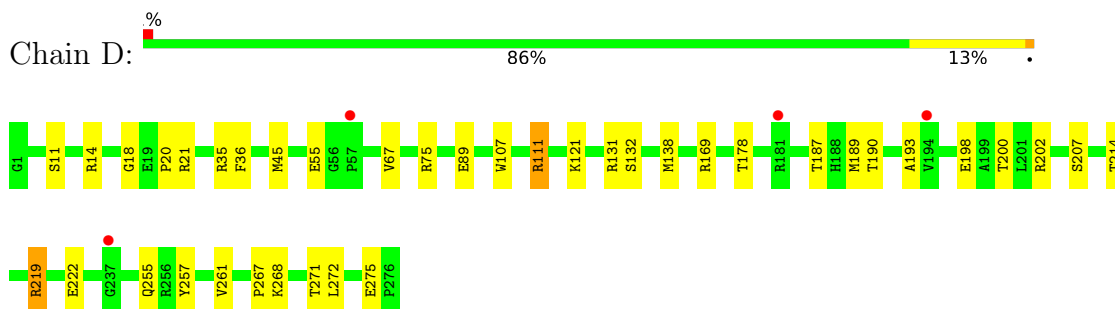
3 Residue-property plots

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.

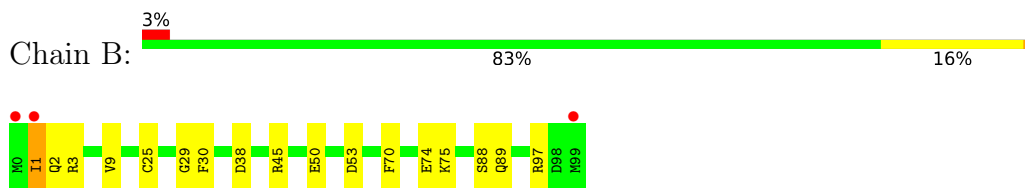
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



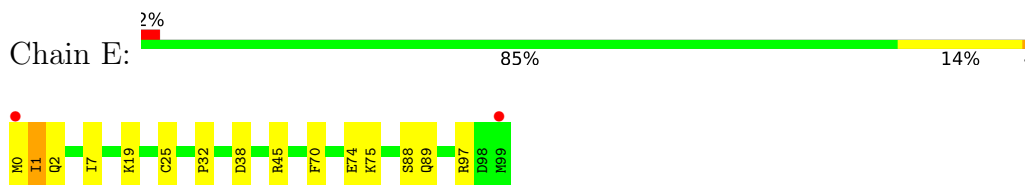
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



- Molecule 3: MET-THR-SER-ALA-ILE-GLY-ILE-LEU-PRO-VAL

Chain C:  90% 10%



• Molecule 3: MET-THR-SER-ALA-ILE-GLY-ILE-LEU-PRO-VAL

Chain F:  70% 20% 10%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	202.94Å 50.28Å 119.07Å 90.00° 122.86° 90.00°	Depositor
Resolution (Å)	41.05 – 2.45 41.05 – 2.45	Depositor EDS
% Data completeness (in resolution range)	98.3 (41.05-2.45) 98.3 (41.05-2.45)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.223 , 0.270 0.227 , 0.268	Depositor DCC
R_{free} test set	1818 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	62.8	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 36.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6354	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.50% 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	1.13	2/2331 (0.1%)	1.16	6/3164 (0.2%)
1	D	1.13	2/2320 (0.1%)	1.15	5/3149 (0.2%)
2	B	1.34	5/860 (0.6%)	1.17	1/1162 (0.1%)
2	E	1.31	3/860 (0.3%)	1.18	1/1162 (0.1%)
3	C	1.21	0/69	1.27	0/92
3	F	1.39	0/69	1.42	1/92 (1.1%)
All	All	1.19	12/6509 (0.2%)	1.16	14/8821 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	2

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	11	SER	CA-C	-7.55	1.43	1.52
2	E	7	ILE	C-O	-6.92	1.17	1.24
2	B	25	CYS	CA-C	6.86	1.61	1.52
1	A	6	ARG	CA-C	-6.26	1.45	1.52
2	E	45	ARG	N-CA	5.86	1.53	1.46
1	A	11	SER	CA-C	-5.51	1.46	1.52
2	E	25	CYS	CA-C	5.48	1.59	1.52
1	D	36	PHE	C-O	-5.44	1.17	1.23
2	B	9	VAL	C-O	-5.43	1.18	1.24
2	B	45	ARG	N-CA	5.21	1.52	1.46
2	B	53	ASP	CB-CG	5.17	1.65	1.52
2	B	30	PHE	C-O	-5.11	1.17	1.23

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	45	ARG	CB-CA-C	-7.74	100.28	109.80
2	E	45	ARG	CB-CA-C	-7.66	100.38	109.80
3	F	6	GLY	N-CA-C	-7.60	98.39	112.37
1	A	224	GLN	CB-CA-C	6.89	120.36	110.79
1	D	138	MET	N-CA-C	6.86	118.84	111.36
1	A	18	GLY	N-CA-C	6.85	121.44	111.24
1	D	18	GLY	N-CA-C	6.79	121.35	111.24
1	D	45	MET	N-CA-C	-5.78	101.06	110.20
1	A	138	MET	N-CA-C	5.74	117.61	111.36
1	D	21	ARG	N-CA-C	-5.48	102.77	110.50
1	D	55	GLU	N-CA-C	5.43	117.90	110.24
1	A	163	THR	N-CA-C	5.37	116.81	111.07
1	A	45	MET	N-CA-C	-5.25	101.90	110.20
1	A	21	ARG	N-CA-C	-5.09	103.32	110.50

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	14	ARG	Peptide
1	D	193	ALA	Peptide

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	2264	0	2109	7	0
1	D	2254	0	2103	7	0
2	B	837	0	803	2	0
2	E	837	0	803	3	0
3	C	69	0	80	0	0
3	F	69	0	80	1	0
4	A	6	0	0	0	0
4	B	4	0	0	0	0
4	C	1	0	0	0	0
4	D	6	0	0	1	0
4	E	7	0	0	0	0
All	All	6354	0	5978	20	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 2.

All (20) 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)
1:A:190:THR:OG1	1:A:202:ARG:HB3	2.09	0.53
1:D:190:THR:OG1	1:D:202:ARG:HB3	2.09	0.52
3:F:7:ILE:O	3:F:7:ILE:HG22	2.12	0.49
1:D:111:ARG:HA	1:D:111:ARG:CZ	2.43	0.48
1:A:74:HIS:HE1	1:A:97:ARG:HE	1.62	0.48
1:A:74:HIS:CE1	1:A:97:ARG:HE	2.32	0.47
1:A:187:THR:HG21	1:A:261:VAL:HG21	1.95	0.47
2:E:0:MET:HE1	2:E:32:PRO:HG3	1.95	0.47
1:D:187:THR:HG21	1:D:261:VAL:HG21	1.98	0.46
1:A:219:ARG:HG2	1:A:257:TYR:CE2	2.51	0.45
1:D:107:TRP:HB3	1:D:169:ARG:HD3	1.99	0.45
1:A:107:TRP:HB3	1:A:169:ARG:HD3	2.00	0.43
2:B:1:ILE:O	2:B:1:ILE:HD13	2.19	0.42
2:E:1:ILE:O	2:E:1:ILE:HD13	2.20	0.42
4:D:305:HOH:O	2:E:0:MET:HG2	2.19	0.42
2:B:3:ARG:HG2	2:B:29:GLY:O	2.20	0.41
1:D:20:PRO:HG2	1:D:75:ARG:HG3	2.03	0.41
1:A:259:CYS:HB3	1:A:272:LEU:HB2	2.03	0.41
1:D:187:THR:HB	1:D:272:LEU:HD21	2.03	0.40
1:D:219:ARG:HG2	1:D:257:TYR:CE2	2.56	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	275/276 (100%)	259 (94%)	14 (5%)	2 (1%)	18 24

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	274/276 (99%)	257 (94%)	15 (6%)	2 (1%)	18	24
2	B	98/100 (98%)	95 (97%)	2 (2%)	1 (1%)	12	14
2	E	98/100 (98%)	95 (97%)	2 (2%)	1 (1%)	12	14
3	C	8/10 (80%)	7 (88%)	1 (12%)	0	100	100
3	F	8/10 (80%)	7 (88%)	0	1 (12%)	0	0
All	All	761/772 (99%)	720 (95%)	34 (4%)	7 (1%)	14	17

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	97	ARG
2	E	97	ARG
1	A	275	GLU
1	D	275	GLU
1	A	267	PRO
1	D	267	PRO
3	F	7	ILE

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	233/232 (100%)	214 (92%)	19 (8%)	10	13
1	D	232/232 (100%)	214 (92%)	18 (8%)	11	14
2	B	95/95 (100%)	86 (90%)	9 (10%)	8	8
2	E	95/95 (100%)	86 (90%)	9 (10%)	8	8
3	C	8/8 (100%)	7 (88%)	1 (12%)	4	4
3	F	8/8 (100%)	7 (88%)	1 (12%)	4	4
All	All	671/670 (100%)	614 (92%)	57 (8%)	10	11

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	VAL
1	A	35	ARG
1	A	67	VAL
1	A	111	ARG
1	A	121	LYS
1	A	131	ARG
1	A	132	SER
1	A	176	LYS
1	A	178	THR
1	A	189	MET
1	A	198	GLU
1	A	200	THR
1	A	207	SER
1	A	214	THR
1	A	219	ARG
1	A	224	GLN
1	A	226	GLN
1	A	255	GLN
1	A	271	THR
2	B	1	ILE
2	B	2	GLN
2	B	38	ASP
2	B	50	GLU
2	B	70	PHE
2	B	74	GLU
2	B	75	LYS
2	B	88	SER
2	B	89	GLN
3	C	5	ILE
1	D	35	ARG
1	D	67	VAL
1	D	89	GLU
1	D	111	ARG
1	D	121	LYS
1	D	131	ARG
1	D	132	SER
1	D	178	THR
1	D	189	MET
1	D	198	GLU
1	D	200	THR
1	D	207	SER
1	D	214	THR
1	D	219	ARG

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Mol	Chain	Res	Type
1	D	222	GLU
1	D	255	GLN
1	D	268	LYS
1	D	271	THR
2	E	1	ILE
2	E	2	GLN
2	E	19	LYS
2	E	38	ASP
2	E	70	PHE
2	E	74	GLU
2	E	75	LYS
2	E	88	SER
2	E	89	GLN
3	F	10	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	74	HIS
1	A	86	ASN
1	A	115	GLN
1	A	155	GLN
1	A	174	ASN
1	A	253	GLN
1	A	255	GLN
2	B	13	HIS
2	B	31	HIS
2	B	83	ASN
1	D	43	GLN
1	D	54	GLN
1	D	74	HIS
1	D	86	ASN
1	D	155	GLN
1	D	174	ASN
1	D	191	HIS
1	D	253	GLN
2	E	31	HIS
2	E	83	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 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	276/276 (100%)	0.09	2 (0%) 84 86	42, 86, 128, 149	1 (0%)
1	D	276/276 (100%)	0.15	4 (1%) 73 75	51, 92, 137, 157	0
2	B	100/100 (100%)	-0.06	3 (3%) 52 53	51, 72, 108, 122	0
2	E	100/100 (100%)	0.00	2 (2%) 65 66	48, 72, 107, 128	0
3	C	10/10 (100%)	0.35	0 100 100	74, 77, 90, 91	0
3	F	10/10 (100%)	0.37	0 100 100	73, 79, 88, 98	0
All	All	772/772 (100%)	0.09	11 (1%) 73 75	42, 84, 130, 157	1 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	188[A]	HIS	2.9
2	B	1	ILE	2.8
1	A	276	PRO	2.7
2	B	99	MET	2.7
1	D	237	GLY	2.6
2	B	0	MET	2.5
2	E	0	MET	2.4
1	D	194	VAL	2.3
2	E	99	MET	2.2
1	D	57	PRO	2.1
1	D	181	ARG	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.