



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:58 AM UTC

PDB ID : 3QUK / pdb_00003quk
Title : Crystal structures of the murine class I major histocompatibility complex H-2Db in complex with LCMV-derived gp33 altered peptide ligand (Y4A)
Authors : Allerbring, E.; Duru, A.D.; Uchtenhagen, H.; Madhurantakam, C.; Grimm, S.; Tomek, M.B.; Mazumdar, P.A.; Spetz, A.; Friemann, R.; Sandalova, T.; Uhlin, M.; Nygren, P.; Achour, A.
Deposited on : 2011-02-24
Resolution : 2.41 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

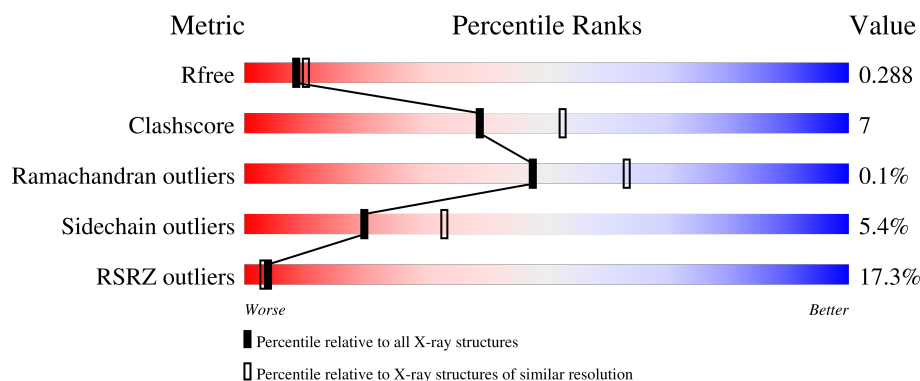
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	17% 80% 13% 5%
1	D	276	19% 79% 14% 5%
2	B	99	18% 74% 24% .
2	E	99	8% 87% 13%
3	C	9	78% 22%

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Mol	Chain	Length	Quality of chain
3	F	9	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6443 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 I histocompatibility antigen, D-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	262	Total	C	N	O	S	0	2	0
			2178	1375	385	409	9			
1	D	261	Total	C	N	O	S	0	2	0
			2173	1373	388	403	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	1	0
			828	528	139	154	7			
2	E	99	Total	C	N	O	S	0	1	0
			832	530	142	153	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	85	ASP	ALA	variant	UNP P01887
E	85	ASP	ALA	variant	UNP P01887

- Molecule 3 is a protein called Pre-glycoprotein polypeptide GP complex.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	S	0	0	0
			66	42	11	12	1			
3	F	9	Total	C	N	O	S	0	0	0
			66	42	11	12	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	4	ALA	TYR	engineered mutation	UNP P07399
C	9	MET	CYS	engineered mutation	UNP P07399

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Chain	Residue	Modelled	Actual	Comment	Reference
F	4	ALA	TYR	engineered mutation	UNP P07399
F	9	MET	CYS	engineered mutation	UNP P07399

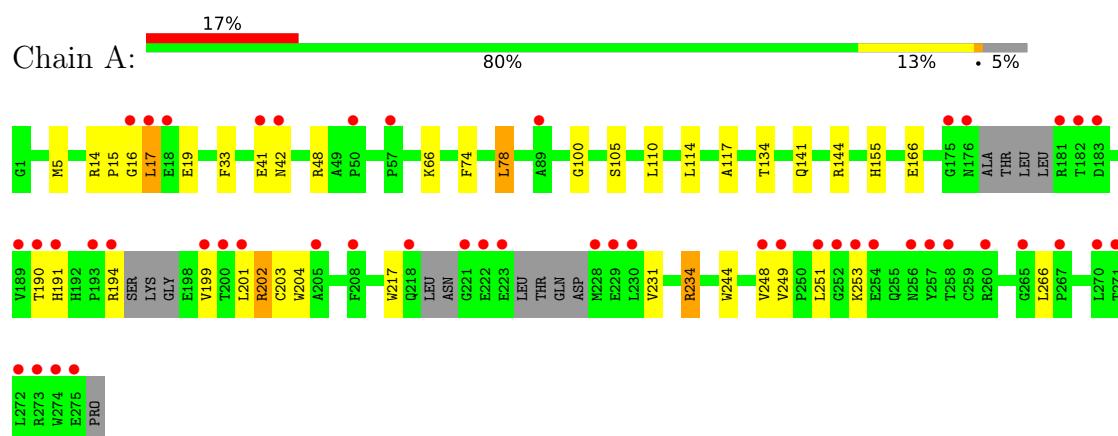
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	110	Total O 110 110	0	0
4	B	44	Total O 44 44	0	0
4	C	8	Total O 8 8	0	0
4	D	87	Total O 87 87	0	0
4	E	49	Total O 49 49	0	0
4	F	2	Total O 2 2	0	0

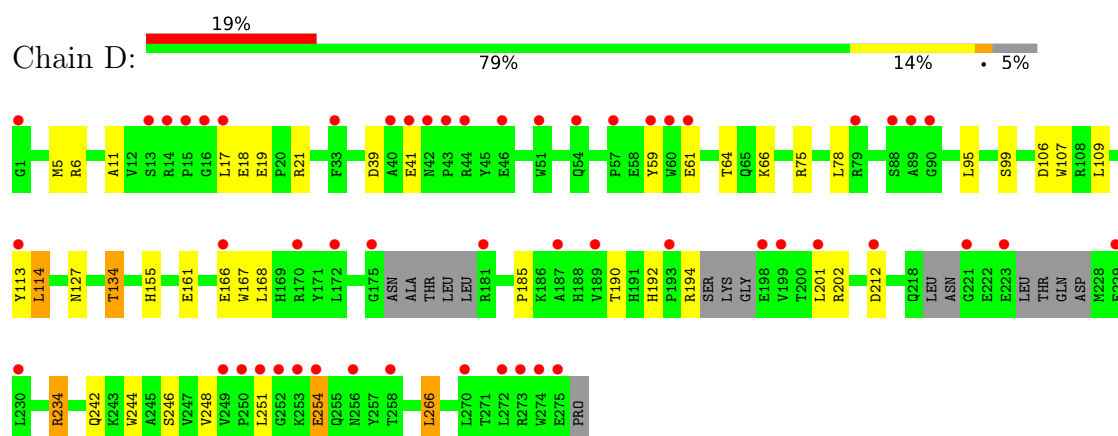
3 Residue-property plots [i](#)

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.

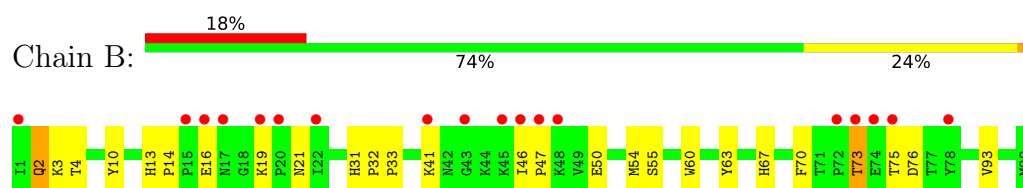
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain



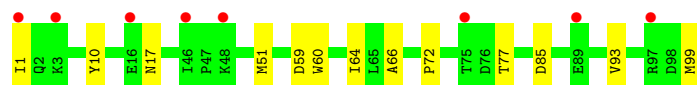
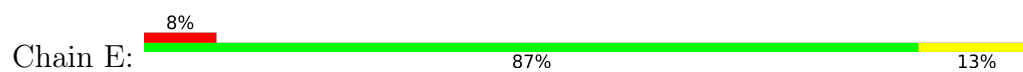
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain



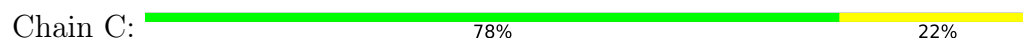
- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



- Molecule 3: Pre-glycoprotein polypeptide GP complex



- Molecule 3: Pre-glycoprotein polypeptide GP complex



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	118.51Å 126.21Å 96.47Å 90.00° 124.54° 90.00°	Depositor
Resolution (Å)	33.90 – 2.41 33.90 – 2.41	Depositor EDS
% Data completeness (in resolution range)	99.0 (33.90-2.41) 98.9 (33.90-2.41)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.242 , 0.285 0.248 , 0.288	Depositor DCC
R_{free} test set	2252 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	43.2	Xtriage
Anisotropy	0.022	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6443	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.73% 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	0/2240	0.76	0/3035
1	D	0.55	0/2238	0.79	0/3030
2	B	0.63	0/854	0.83	1/1158 (0.1%)
2	E	0.57	0/858	0.83	0/1162
3	C	0.72	0/66	0.82	0/86
3	F	0.85	0/66	0.79	0/86
All	All	0.58	0/6322	0.79	1/8557 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	41	LYS	CB-CA-C	-6.96	99.53	110.74

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	2178	0	2034	28	0
1	D	2173	0	2042	37	0
2	B	828	0	802	15	0
2	E	832	0	808	5	0
3	C	66	0	70	2	0
3	F	66	0	70	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	110	0	0	3	0
4	B	44	0	0	2	0
4	C	8	0	0	0	0
4	D	87	0	0	0	0
4	E	49	0	0	0	0
4	F	2	0	0	0	0
All	All	6443	0	5826	81	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 7.

All (81) 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:D:19:GLU:CG	1:D:75[B]:ARG:HH21	1.31	1.42
1:D:19:GLU:CG	1:D:75[B]:ARG:NH2	2.00	1.23
1:D:19:GLU:HG3	1:D:75[B]:ARG:NH2	1.65	1.08
1:D:19:GLU:HG2	1:D:75[B]:ARG:HH21	0.93	1.07
1:D:19:GLU:HG2	1:D:75[B]:ARG:NH2	1.66	1.02
1:A:144:ARG:HD2	4:A:387:HOH:O	1.64	0.94
1:A:14:ARG:HB2	1:A:17:LEU:HD22	1.61	0.82
1:D:59:TYR:CE2	3:F:1:LYS:HE3	2.19	0.77
1:A:41:GLU:HG2	1:A:42:ASN:N	2.01	0.73
1:D:59:TYR:CD2	3:F:1:LYS:HE3	2.26	0.71
1:D:167:TRP:CE2	3:F:1:LYS:HD3	2.27	0.69
1:D:41:GLU:CD	1:D:41:GLU:H	2.02	0.68
1:D:155:HIS:HB3	3:F:6:PHE:CZ	2.29	0.68
1:D:127:ASN:OD1	1:D:134:THR:HG22	1.95	0.67
1:A:41:GLU:HG2	1:A:42:ASN:H	1.60	0.66
2:B:3:LYS:HE3	4:B:213:HOH:O	1.95	0.66
1:A:16:GLY:O	1:A:17:LEU:C	2.38	0.65
2:B:73:THR:HB	2:B:76:ASP:OD2	1.97	0.65
1:D:18:GLU:O	1:D:18:GLU:HG2	1.97	0.64
1:A:202:ARG:HD2	1:A:244:TRP:CD2	2.35	0.62
1:D:59:TYR:CE2	3:F:1:LYS:CE	2.82	0.62
1:D:19:GLU:HG3	1:D:75[B]:ARG:CZ	2.30	0.61
1:D:99:SER:HB3	1:D:114:LEU:HD23	1.80	0.61
1:D:202:ARG:HD3	1:D:246:SER:HB3	1.84	0.60
1:A:202:ARG:HG2	1:A:204:TRP:NE1	2.17	0.60
1:A:19:GLU:HA	1:A:19:GLU:OE1	2.02	0.60
1:A:14:ARG:HB2	1:A:17:LEU:HD13	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:19:GLU:CD	1:D:75[B]:ARG:HE	2.12	0.57
1:D:5:MET:HB2	1:D:168:LEU:HD13	1.86	0.57
1:A:201:LEU:HD12	1:A:249:VAL:HG21	1.86	0.56
1:A:14:ARG:HB2	1:A:17:LEU:CD2	2.35	0.55
1:D:201:LEU:HD11	1:D:254:GLU:HB2	1.90	0.54
1:D:234:ARG:HD3	2:E:10:TYR:CE2	2.42	0.54
1:A:155:HIS:HB3	3:C:6:PHE:CZ	2.43	0.53
2:B:73:THR:CB	2:B:76:ASP:OD2	2.58	0.52
2:B:33:PRO:O	2:B:54:MET:HE1	2.11	0.51
1:D:202:ARG:HD2	1:D:244:TRP:CE3	2.46	0.51
1:D:202:ARG:HD2	1:D:244:TRP:CD2	2.46	0.50
1:A:144:ARG:CD	4:A:387:HOH:O	2.38	0.50
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.46	0.50
1:D:167:TRP:CZ2	3:F:1:LYS:HD3	2.47	0.50
1:A:194:ARG:HH11	1:A:248:VAL:HG21	1.76	0.49
4:A:397:HOH:O	1:D:212:ASP:HB3	2.11	0.49
1:D:11:ALA:HA	1:D:21:ARG:O	2.13	0.49
1:A:191:HIS:CE1	1:A:199:VAL:HG21	2.47	0.49
1:A:66:LYS:HD2	3:C:4:ALA:HB2	1.94	0.48
1:A:33:PHE:C	1:A:48:ARG:HB2	2.38	0.48
1:A:41:GLU:CG	1:A:42:ASN:H	2.26	0.48
1:D:234:ARG:HD2	1:D:242:GLN:HB2	1.97	0.46
1:A:41:GLU:CG	1:A:42:ASN:N	2.74	0.46
2:B:50:GLU:HB2	2:B:67:HIS:CE1	2.50	0.46
2:B:55:SER:HB2	2:B:63:TYR:CZ	2.51	0.45
2:B:16:GLU:OE1	2:B:19:LYS:HD2	2.16	0.45
1:D:19:GLU:CD	1:D:75[B]:ARG:HH21	2.10	0.44
1:A:251:LEU:C	1:A:253:LYS:H	2.24	0.44
2:E:17:ASN:HA	2:E:72:PRO:O	2.18	0.44
2:E:59:ASP:O	2:E:60:TRP:HB2	2.18	0.44
1:A:141:GLN:OE1	1:A:144:ARG:NH1	2.52	0.43
1:A:5:MET:O	1:A:100:GLY:HA3	2.18	0.43
1:D:234:ARG:NH2	2:E:99:MET:O	2.41	0.43
1:A:190:THR:OG1	1:A:202:ARG:HB3	2.19	0.42
1:D:190:THR:OG1	1:D:192:HIS:HE1	2.02	0.42
1:A:194:ARG:NH1	1:A:248:VAL:HG21	2.33	0.42
1:D:194:ARG:HE	1:D:248:VAL:HG11	1.84	0.42
2:B:2:GLN:H	2:B:2:GLN:HG2	1.53	0.42
1:D:59:TYR:HE2	3:F:1:LYS:HE3	1.81	0.42
1:D:59:TYR:CD2	3:F:1:LYS:CE	3.01	0.42
2:B:32:PRO:HB2	2:B:33:PRO:HD2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:106:ASP:O	1:D:107:TRP:HB2	2.20	0.42
1:D:95:LEU:HD11	3:F:9:MET:HE3	2.02	0.41
1:D:185:PRO:HD2	1:D:266:LEU:HD13	2.01	0.41
1:A:234:ARG:HD3	2:B:10:TYR:CE2	2.56	0.41
1:A:203:CYS:HB2	1:A:217:TRP:CZ2	2.56	0.41
2:B:13:HIS:O	2:B:14:PRO:C	2.64	0.41
1:D:109:LEU:HD22	1:D:161:GLU:HG2	2.02	0.41
2:E:51:MET:SD	2:E:66:ALA:HB2	2.61	0.41
2:B:46:ILE:HA	2:B:47:PRO:HD3	1.85	0.40
2:B:21:ASN:HB3	2:B:70:PHE:CE1	2.55	0.40
1:D:6:ARG:NH2	1:D:113:TYR:CE1	2.89	0.40
1:A:74:PHE:O	1:A:78:LEU:HB2	2.20	0.40
2:B:31:HIS:ND1	4:B:116:HOH:O	2.37	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	254/276 (92%)	244 (96%)	9 (4%)	1 (0%)	30	42
1	D	253/276 (92%)	245 (97%)	8 (3%)	0	100	100
2	B	98/99 (99%)	94 (96%)	4 (4%)	0	100	100
2	E	98/99 (99%)	95 (97%)	3 (3%)	0	100	100
3	C	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	F	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
All	All	717/768 (93%)	690 (96%)	26 (4%)	1 (0%)	48	63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	17	LEU

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	224/234 (96%)	213 (95%)	11 (5%)	22	37
1	D	223/234 (95%)	210 (94%)	13 (6%)	18	30
2	B	95/94 (101%)	90 (95%)	5 (5%)	20	34
2	E	95/94 (101%)	90 (95%)	5 (5%)	20	34
3	C	6/6 (100%)	6 (100%)	0	100	100
3	F	6/6 (100%)	5 (83%)	1 (17%)	2	2
All	All	649/668 (97%)	614 (95%)	35 (5%)	20	33

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	PRO
1	A	78	LEU
1	A	105	SER
1	A	110	LEU
1	A	114	LEU
1	A	134	THR
1	A	166	GLU
1	A	202	ARG
1	A	231	VAL
1	A	234	ARG
1	A	266	LEU
2	B	2	GLN
2	B	4	THR
2	B	73	THR
2	B	75	THR
2	B	93	VAL
1	D	17	LEU
1	D	39	ASP

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Mol	Chain	Res	Type
1	D	61	GLU
1	D	64	THR
1	D	66	LYS
1	D	78	LEU
1	D	114	LEU
1	D	134	THR
1	D	166	GLU
1	D	234	ARG
1	D	251	LEU
1	D	254	GLU
1	D	266	LEU
2	E	1	ILE
2	E	64	ILE
2	E	77	THR
2	E	85	ASP
2	E	93	VAL
3	F	1	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	97	GLN
1	A	191	HIS
1	A	192	HIS
2	B	31	HIS
3	C	5	ASN
1	D	97	GLN
1	D	192	HIS
1	D	256	ASN
2	E	13	HIS
2	E	34	HIS
2	E	67	HIS
3	F	5	ASN

5.3.3 RNA ⓘ

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 ⓘ

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	262/276 (94%)	1.10	48 (18%)	3 3	19, 52, 84, 117	2 (0%)
1	D	261/276 (94%)	1.25	53 (20%)	3 2	20, 56, 82, 105	2 (0%)
2	B	99/99 (100%)	1.19	18 (18%)	3 3	27, 54, 66, 68	1 (1%)
2	E	99/99 (100%)	0.99	8 (8%)	18 14	27, 54, 65, 69	1 (1%)
3	C	9/9 (100%)	0.55	0	100 100	42, 47, 49, 49	0
3	F	9/9 (100%)	0.87	1 (11%)	10 8	45, 49, 55, 58	0
All	All	739/768 (96%)	1.14	128 (17%)	4 3	19, 54, 81, 117	6 (0%)

All (128) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	221	GLY	7.7
1	A	221	GLY	6.9
1	A	41	GLU	6.1
1	A	17	LEU	5.1
1	D	41	GLU	5.0
1	D	90	GLY	4.9
1	D	251	LEU	4.7
2	B	73	THR	4.6
1	A	223	GLU	4.6
1	A	272	LEU	4.5
1	D	274	TRP	4.3
2	B	1	ILE	4.3
2	E	1	ILE	4.3
1	D	89	ALA	4.3
1	D	79[A]	ARG	4.1
1	D	275	GLU	4.0
2	E	97[A]	ARG	4.0
1	A	257	TYR	3.9
1	D	40	ALA	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	193	PRO	3.9
1	D	175	GLY	3.8
1	A	274	TRP	3.8
1	D	16	GLY	3.8
1	D	249	VAL	3.6
1	A	229[A]	GLU	3.6
1	A	265	GLY	3.5
1	A	249	VAL	3.5
1	D	250	PRO	3.5
1	D	51	TRP	3.4
1	A	230	LEU	3.3
1	A	89	ALA	3.3
1	A	253	LYS	3.3
1	D	256	ASN	3.2
1	A	267	PRO	3.2
2	B	47	PRO	3.2
1	D	54	GLN	3.1
1	D	59	TYR	3.0
1	A	251	LEU	3.0
1	D	13	SER	3.0
1	A	176	ASN	3.0
1	D	193	PRO	3.0
2	B	19	LYS	2.9
1	A	181	ARG	2.9
1	D	187	ALA	2.9
1	A	190	THR	2.9
2	B	48	LYS	2.9
1	D	230	LEU	2.9
2	B	16	GLU	2.9
1	D	61	GLU	2.9
1	D	223	GLU	2.9
1	A	16	GLY	2.8
1	A	273	ARG	2.8
1	A	191	HIS	2.8
2	B	41	LYS	2.8
1	D	88	SER	2.8
1	A	189	VAL	2.8
1	A	270	LEU	2.8
1	D	17	LEU	2.8
1	A	254	GLU	2.8
1	D	15	PRO	2.7
1	A	271	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	253	LYS	2.7
1	D	252	GLY	2.7
1	A	194	ARG	2.7
2	B	75	THR	2.7
1	D	272	LEU	2.7
1	A	50	PRO	2.6
1	A	222	GLU	2.6
1	A	182	THR	2.6
1	D	254	GLU	2.6
1	A	252	GLY	2.6
1	D	181	ARG	2.6
1	A	228	MET	2.6
1	A	258	THR	2.5
2	B	22	ILE	2.5
1	A	275	GLU	2.5
1	D	166	GLU	2.5
1	D	14	ARG	2.4
1	A	200	THR	2.4
2	E	75	THR	2.4
1	D	198	GLU	2.4
1	D	113	TYR	2.4
1	D	44	ARG	2.4
1	D	1	GLY	2.4
1	D	57	PRO	2.4
2	B	74	GLU	2.4
1	A	208	PHE	2.4
1	D	43	PRO	2.3
2	E	48	LYS	2.3
3	F	1	LYS	2.3
1	D	273	ARG	2.3
1	A	218	GLN	2.3
1	D	189	VAL	2.3
1	A	42	ASN	2.3
1	D	42	ASN	2.3
1	A	260	ARG	2.3
1	A	175	GLY	2.3
1	A	57	PRO	2.3
1	A	201	LEU	2.2
1	D	46	GLU	2.2
2	B	20	PRO	2.2
1	A	256	ASN	2.2
1	A	18	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
2	E	16	GLU	2.2
2	E	89	GLU	2.2
2	B	45	LYS	2.2
2	B	17	ASN	2.1
1	D	201	LEU	2.1
1	D	258	THR	2.1
1	D	229	GLU	2.1
1	D	60	TRP	2.1
1	A	183	ASP	2.1
2	B	15	PRO	2.1
1	D	199	VAL	2.1
1	D	172	LEU	2.1
1	D	270	LEU	2.1
2	E	46	ILE	2.1
2	B	78	TYR	2.1
2	E	3	LYS	2.1
1	A	199	VAL	2.1
1	A	248	VAL	2.1
1	D	33	PHE	2.1
2	B	72	PRO	2.1
1	A	205	ALA	2.0
2	B	43	GLY	2.0
2	B	46	ILE	2.0
1	D	212	ASP	2.0
1	D	170	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.