



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 2, 2023 – 11:45 PM EDT

PDB ID : 3QUK
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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.35
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) : 2.35

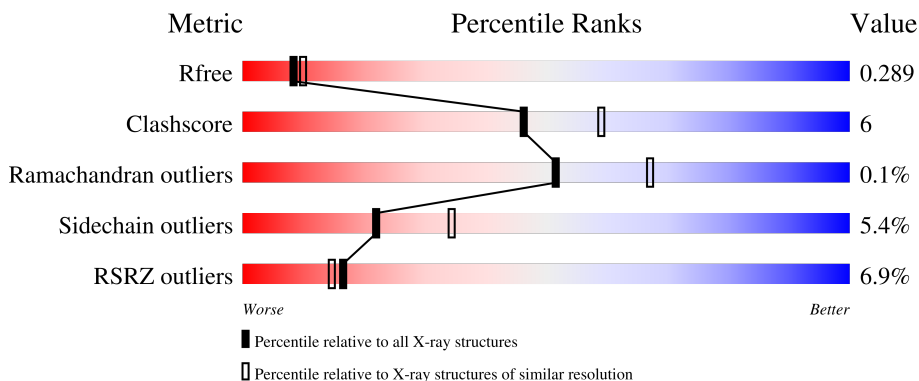
1 Overall quality at a glance i

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.



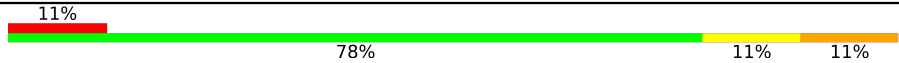
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4647 (2.44-2.40)
Clashscore	141614	5161 (2.44-2.40)
Ramachandran outliers	138981	5073 (2.44-2.40)
Sidechain outliers	138945	5074 (2.44-2.40)
RSRZ outliers	127900	4543 (2.44-2.40)

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	 7% 81% 13% • 5%
1	D	276	 9% 79% 14% • 5%
2	B	99	 5% 77% 20% •
2	E	99	 2% 89% 11%
3	C	9	 78% 22%

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Mol	Chain	Length	Quality of chain
3	F	9	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into four segments: a red segment on the left labeled '11%', a large green segment labeled '78%', a yellow segment labeled '11%', and an orange segment on the right labeled '11%'.</p>

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
			Total	C	N	O	S			
1	A	262	Total 2178	C 1375	N 385	O 409	S 9	0	2	0
1	D	261	Total 2173	C 1373	N 388	O 403	S 9	0	2	0

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

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

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
			Total	C	N	O	S			
3	C	9	Total 66	C 42	N 11	O 12	S 1	0	0	0
3	F	9	Total 66	C 42	N 11	O 12	S 1	0	0	0

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

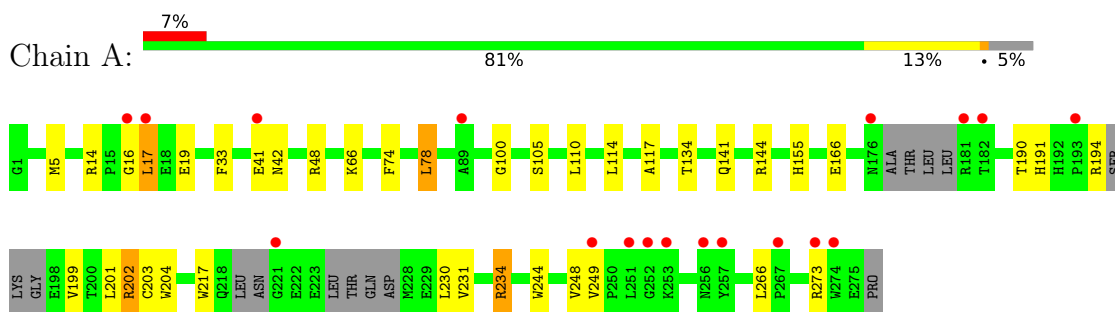
- Molecule 4 is water.

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

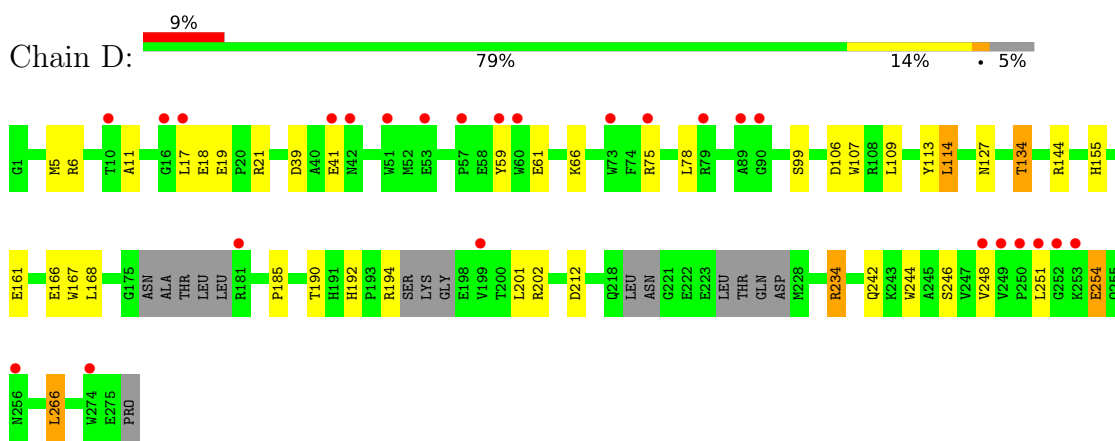
3 Residue-property plots [i](#)

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.

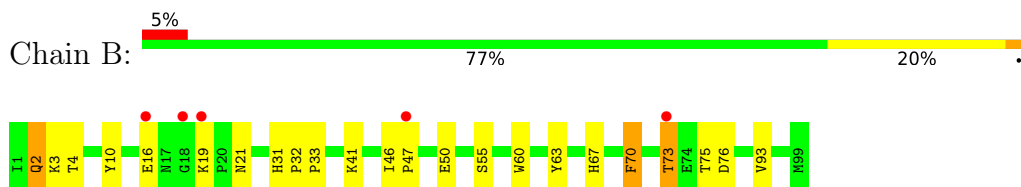
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain



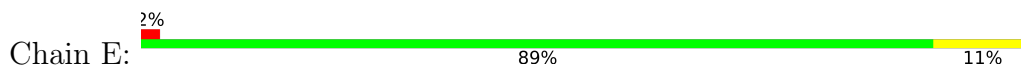
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain



- Molecule 2: Beta-2-microglobulin

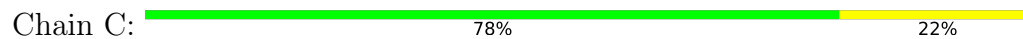


- Molecule 2: Beta-2-microglobulin

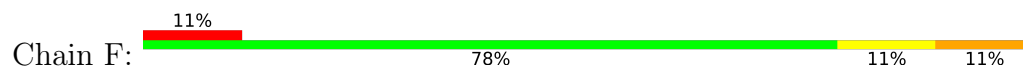




- Molecule 3: Pre-glycoprotein polyprotein GP complex



- Molecule 3: Pre-glycoprotein polyprotein 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 33.89 – 2.41	Depositor EDS
% Data completeness (in resolution range)	99.0 (33.90-2.41) 99.0 (33.89-2.41)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 2.42Å)	Xtrriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.242 , 0.285 0.248 , 0.289	Depositor DCC
R_{free} test set	2252 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	43.2	Xtrriage
Anisotropy	0.022	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
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

Xtrriage'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	0.48	0/2240	0.59	0/3035
1	D	0.47	0/2238	0.61	0/3030
2	B	0.53	0/854	0.63	1/1158 (0.1%)
2	E	0.49	0/858	0.60	0/1162
3	C	0.68	0/66	0.65	0/86
3	F	0.77	0/66	0.72	0/86
All	All	0.49	0/6322	0.61	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	-5.43	99.53	110.40

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	27	0
1	D	2173	0	2042	35	0
2	B	828	0	802	13	0
2	E	832	0	808	5	0
3	C	66	0	70	2	0
3	F	66	0	70	8	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	76	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 (76) 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: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
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:D:41:GLU:CD	1:D:41:GLU:H	2.02	0.63
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:A:16:GLY:O	1:A:17:LEU:C	2.38	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: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
1:D:19:GLU:CD	1:D:75[B]:ARG:HE	2.12	0.53
2:B:73:THR:CB	2:B:76:ASP:OD2	2.58	0.52
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: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
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:33:PHE:C	1:A:48:ARG:HB2	2.38	0.43
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
1:D:106:ASP:O	1:D:107:TRP:HB2	2.20	0.42
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:31:HIS:ND1	4:B:116:HOH:O	2.37	0.41
1:A:203:CYS:HB2	1:A:217:TRP:CZ2	2.56	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

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%)	34	47
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%)	51	67

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	17	LEU

5.3.2 Protein sidechains [i](#)

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%)	212 (95%)	12 (5%)	22	34
1	D	223/234 (95%)	210 (94%)	13 (6%)	20	31
2	B	95/94 (101%)	89 (94%)	6 (6%)	18	27
2	E	95/94 (101%)	92 (97%)	3 (3%)	39	57
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%)	22	34

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

Mol	Chain	Res	Type
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	230	LEU
1	A	231	VAL
1	A	234	ARG
1	A	266	LEU
1	A	273	ARG
2	B	2	GLN
2	B	4	THR
2	B	70	PHE
2	B	73	THR
2	B	75	THR
2	B	93	VAL
1	D	17	LEU
1	D	39	ASP
1	D	61	GLU
1	D	66	LYS

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Mol	Chain	Res	Type
1	D	78	LEU
1	D	114	LEU
1	D	134	THR
1	D	144	ARG
1	D	166	GLU
1	D	234	ARG
1	D	251	LEU
1	D	254	GLU
1	D	266	LEU
2	E	64	ILE
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 (11) 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
1	A	218	GLN
2	B	31	HIS
3	C	5	ASN
1	D	97	GLN
1	D	192	HIS
2	E	38	GLN
3	F	5	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](#)

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%)	0.39	18 (6%) 16 15	24, 52, 84, 117	0
1	D	261/276 (94%)	0.54	25 (9%) 8 7	20, 56, 82, 105	0
2	B	99/99 (100%)	0.28	5 (5%) 28 26	38, 54, 66, 68	0
2	E	99/99 (100%)	0.22	2 (2%) 65 62	40, 54, 65, 69	0
3	C	9/9 (100%)	0.66	0 100 100	42, 47, 49, 49	0
3	F	9/9 (100%)	1.19	1 (11%) 5 4	45, 49, 55, 58	0
All	All	739/768 (96%)	0.42	51 (6%) 16 15	20, 54, 81, 117	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	90	GLY	5.0
1	A	17	LEU	4.4
1	A	251	LEU	4.1
1	A	181	ARG	4.0
1	A	274	TRP	3.9
1	D	256	ASN	3.9
1	D	251	LEU	3.9
1	A	41	GLU	3.9
1	D	51	TRP	3.4
1	D	59	TYR	3.4
1	D	57	PRO	3.3
1	A	89	ALA	3.3
1	D	89	ALA	3.2
1	A	182	THR	3.2
1	D	274	TRP	3.2
1	D	253	LYS	3.1
1	D	181	ARG	3.1
1	D	250	PRO	3.0
1	D	53	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	253	LYS	3.0
1	D	199	VAL	2.9
1	A	257	TYR	2.9
1	D	41	GLU	2.7
1	D	249	VAL	2.7
2	E	48	LYS	2.7
1	A	176	ASN	2.6
2	B	73	THR	2.6
1	D	60	TRP	2.6
1	D	16	GLY	2.6
1	D	42	ASN	2.5
1	A	252	GLY	2.5
2	B	47	PRO	2.4
2	B	16	GLU	2.4
1	A	193	PRO	2.4
1	A	256	ASN	2.3
1	A	249	VAL	2.3
1	D	248	VAL	2.3
1	D	75[A]	ARG	2.3
1	A	221	GLY	2.3
1	A	273	ARG	2.2
1	D	252	GLY	2.2
1	A	16	GLY	2.1
1	D	73	TRP	2.1
3	F	5	ASN	2.1
2	E	47	PRO	2.1
1	D	17	LEU	2.1
1	D	79[A]	ARG	2.1
1	D	10	THR	2.1
2	B	18	GLY	2.0
2	B	19	LYS	2.0
1	A	267	PRO	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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.