



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 08:59 PM UTC

PDB ID : 5TIL / pdb_00005til
Title : Murine class I major histocompatibility complex H-2 Db in complex with LCMV-derived GP33 altered peptide V3P and T-cell receptor P14
Authors : Achour, A.; Sandalova, T.; Allerbring, E.B.
Deposited on : 2016-10-03
Resolution : 2.83 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

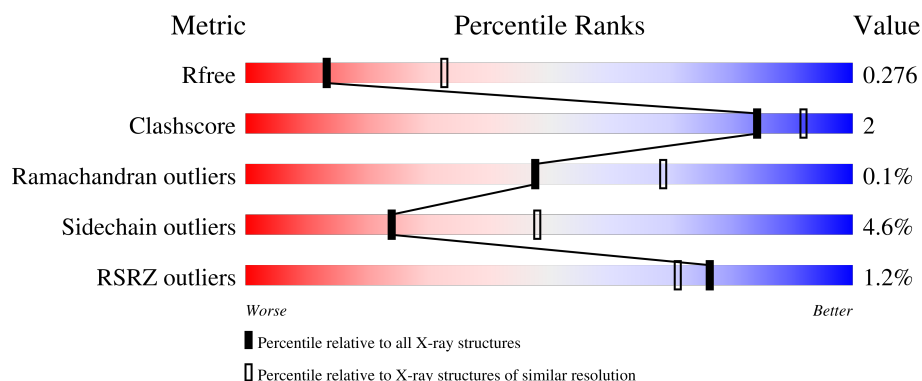
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.83 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	1520 (2.86-2.82)
Clashscore	190562	1559 (2.86-2.82)
Ramachandran outliers	187476	1517 (2.86-2.82)
Sidechain outliers	187428	1518 (2.86-2.82)
RSRZ outliers	180081	1521 (2.86-2.82)

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	2% 90% 9%
1	D	276	% 89% 10% .
2	B	99	93% 6% .
2	E	99	93% 6% .
3	C	9	78% 11% 11%

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Mol	Chain	Length	Quality of chain
3	F	9	89% 11%
4	G	205	47% 8% 44%
5	H	238	3% 87% 10%
5	L	238	% 89% 9%
6	K	160	% 61% 12% 26%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11851 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	275	Total	C	N	O	S	0	0	0
			2260	1428	399	424	9			
1	D	275	Total	C	N	O	S	0	0	0
			2260	1428	399	424	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	98	Total	C	N	O	S	0	0	0
			813	518	137	151	7			
2	E	98	Total	C	N	O	S	0	0	0
			809	515	136	151	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
			73	48	11	13	1			
3	F	9	Total	C	N	O	S	0	0	0
			73	48	11	13	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	3	PRO	VAL	engineered mutation	UNP Q9QDK7
C	9	MET	CYS	engineered mutation	UNP Q9QDK7

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Chain	Residue	Modelled	Actual	Comment	Reference
F	3	PRO	VAL	engineered mutation	UNP Q9QDK7
F	9	MET	CYS	engineered mutation	UNP Q9QDK7

- Molecule 4 is a protein called alpha chain of P14 T cell receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	G	114	Total	C	N	O	S	0	0	0
			905	580	147	176	2			

- Molecule 5 is a protein called Beta chain of murine T cell receptor P14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	H	230	Total	C	N	O	S	0	0	0
			1814	1139	324	345	6			
5	L	236	Total	C	N	O	S	0	0	0
			1866	1173	332	355	6			

- Molecule 6 is a protein called Alpha chain of murine P14 T cell receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	K	118	Total	C	N	O	S	0	0	0
			938	600	154	182	2			

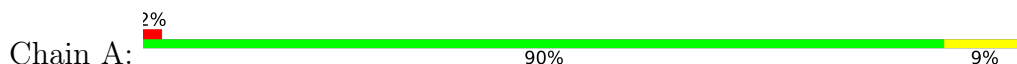
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	9	Total	O	0	0
			9	9		
7	B	5	Total	O	0	0
			5	5		
7	D	5	Total	O	0	0
			5	5		
7	E	7	Total	O	0	0
			7	7		
7	G	2	Total	O	0	0
			2	2		
7	H	5	Total	O	0	0
			5	5		
7	K	5	Total	O	0	0
			5	5		
7	L	2	Total	O	0	0
			2	2		

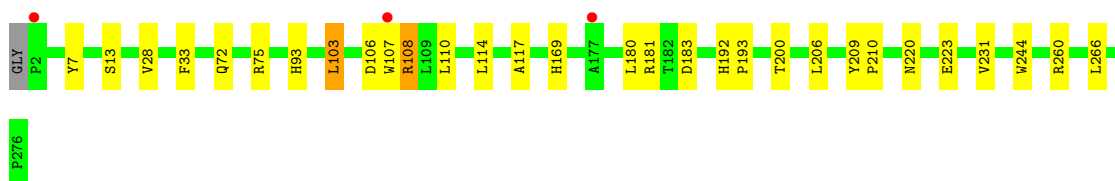
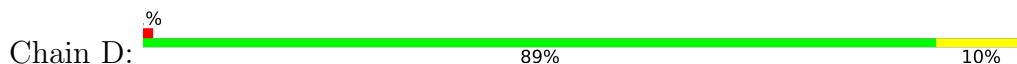
3 Residue-property plots [i](#)

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



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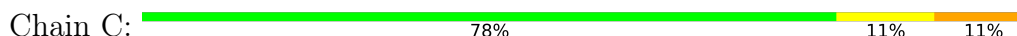
- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



- Molecule 3: Pre-glycoprotein polypeptide GP complex

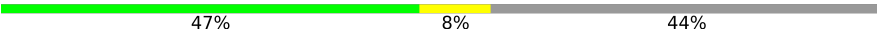


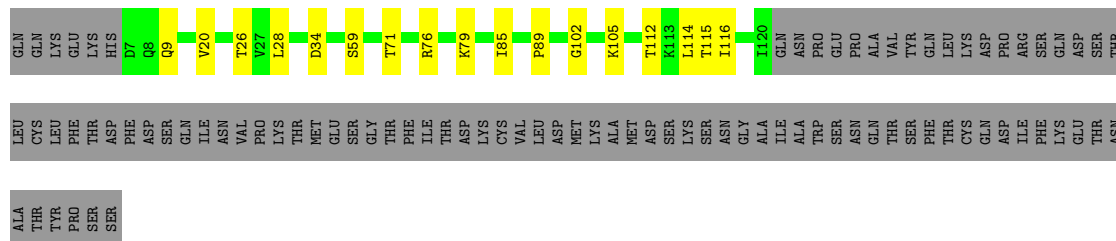
- Molecule 3: Pre-glycoprotein polypeptide GP complex

Chain F:  89% 11%



- Molecule 4: alpha chain of P14 T cell receptor

Chain G:  47% 8% 44%



- Molecule 5: Beta chain of murine T cell receptor P14

Chain H:  3% 87% 10% .



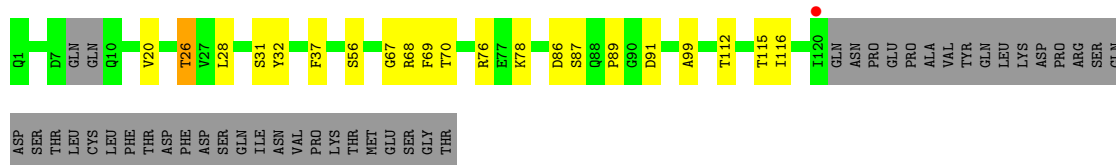
- Molecule 5: Beta chain of murine T cell receptor P14

Chain L:  0% 89% 9% .



- Molecule 6: Alpha chain of murine P14 T cell receptor

Chain K:  0% 61% 12% 26%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.22Å 66.70Å 523.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	52.49 – 2.83 52.49 – 2.83	Depositor EDS
% Data completeness (in resolution range)	97.5 (52.49-2.83) 97.5 (52.49-2.83)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.11 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0124	Depositor
R, R_{free}	0.219 , 0.275 0.222 , 0.276	Depositor DCC
R_{free} test set	2615 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	69.1	Xtriage
Anisotropy	0.097	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 55.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11851	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.60% 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

5.1 Standard geometry

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	0.69	0/2327	0.89	0/3160
1	D	0.68	0/2327	0.86	0/3160
2	B	0.76	0/839	0.92	0/1137
2	E	0.74	0/835	0.97	3/1133 (0.3%)
3	C	0.62	0/75	0.95	0/99
3	F	0.71	0/75	0.92	0/99
4	G	0.73	0/928	0.90	0/1258
5	H	0.70	0/1861	0.89	3/2525 (0.1%)
5	L	0.71	0/1917	0.86	3/2604 (0.1%)
6	K	0.76	0/961	0.91	1/1300 (0.1%)
All	All	0.71	0/12145	0.89	10/16475 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	H	134	ASN	N-CA-C	11.32	123.60	111.03
6	K	86	ASP	N-CA-C	-6.72	102.14	110.41
5	H	158	VAL	N-CA-C	6.44	122.73	109.34
2	E	44	LYS	CB-CA-C	6.11	119.91	110.14
2	E	85	ASP	N-CA-C	5.63	118.14	111.33
2	E	44	LYS	N-CA-C	-5.55	99.25	108.34
5	L	5	SER	CA-C-N	-5.05	113.52	119.84
5	L	5	SER	C-N-CA	-5.05	113.52	119.84
5	L	235	ARG	N-CA-C	5.04	116.25	107.93
5	H	152	VAL	N-CA-C	5.04	115.11	107.75

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	2260	0	2131	12	0
1	D	2260	0	2131	13	0
2	B	813	0	782	2	0
2	E	809	0	771	1	0
3	C	73	0	72	3	0
3	F	73	0	72	1	0
4	G	905	0	875	5	0
5	H	1814	0	1729	7	0
5	L	1866	0	1778	7	0
6	K	938	0	905	10	0
7	A	9	0	0	1	0
7	B	5	0	0	0	0
7	D	5	0	0	0	0
7	E	7	0	0	0	0
7	G	2	0	0	1	0
7	H	5	0	0	0	0
7	K	5	0	0	0	0
7	L	2	0	0	0	0
All	All	11851	0	11246	54	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 (54) 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)
6:K:87:SER:O	6:K:116:ILE:HG21	1.90	0.72
4:G:26:THR:HG21	4:G:114:LEU:HD11	1.71	0.71
6:K:89:PRO:HA	6:K:116:ILE:HB	1.79	0.64
1:A:121:ARG:NH2	7:A:301:HOH:O	2.32	0.62
6:K:68:ARG:NH2	6:K:91:ASP:OD2	2.33	0.60
1:A:72:GLN:OE1	1:A:75:ARG:NH1	2.34	0.60
1:D:107:TRP:HB3	1:D:169:HIS:CE1	2.37	0.59
1:D:72:GLN:OE1	1:D:75:ARG:NH1	2.36	0.58
1:A:107:TRP:HB3	1:A:169:HIS:CE1	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:158:VAL:O	5:H:200:HIS:HB3	2.04	0.57
1:A:123:TYR:CD2	3:C:9:MET:HE2	2.41	0.55
6:K:20:VAL:HG21	6:K:26:THR:HG22	1.88	0.55
6:K:68:ARG:NH2	6:K:91:ASP:OD1	2.42	0.53
5:L:68:SER:HB3	5:L:71:ASN:HD22	1.75	0.51
6:K:68:ARG:NH2	6:K:91:ASP:CG	2.70	0.49
1:A:192:HIS:O	1:A:193:PRO:C	2.55	0.49
5:H:107:ARG:NH1	5:H:150:ASP:O	2.45	0.49
6:K:37:PHE:CD2	6:K:99:ALA:HB1	2.47	0.48
5:H:58:ILE:HG22	5:H:58:ILE:O	2.13	0.48
1:D:28:VAL:HG23	1:D:33:PHE:CE1	2.49	0.48
5:L:19:LEU:HD22	5:L:106:THR:HG21	1.94	0.48
4:G:28:LEU:HD12	4:G:112:THR:HG21	1.95	0.47
1:D:220:ASN:HB3	7:G:301:HOH:O	2.14	0.47
1:D:117:ALA:HB2	2:E:60:TRP:CE2	2.50	0.47
5:L:116:ASN:HD22	5:L:116:ASN:N	2.13	0.47
1:D:181:ARG:NH1	1:D:183:ASP:OD2	2.49	0.46
5:L:144:ALA:HB2	5:L:205:VAL:HG21	1.98	0.45
1:D:106:ASP:O	1:D:108:ARG:N	2.47	0.45
4:G:89:PRO:HA	4:G:116:ILE:HB	1.99	0.44
1:D:7:TYR:CE1	3:F:2:ALA:HB2	2.53	0.44
1:A:28:VAL:HG23	1:A:33:PHE:CE1	2.53	0.44
4:G:34:ASP:C	4:G:34:ASP:OD1	2.61	0.44
1:D:13:SER:HB2	1:D:93:HIS:H	1.82	0.44
1:A:209:TYR:CD1	1:A:210:PRO:HA	2.53	0.44
6:K:28:LEU:HD22	6:K:112:THR:HG21	1.99	0.43
1:D:209:TYR:CD1	1:D:210:PRO:HA	2.53	0.43
1:A:106:ASP:O	1:A:108:ARG:N	2.49	0.43
5:H:62:TYR:HB3	5:H:74:LEU:HD11	2.00	0.43
5:L:34:ARG:HB2	5:L:44:ILE:HD11	2.01	0.43
5:H:158:VAL:HG23	5:H:159:ASN:H	1.83	0.42
1:A:233:THR:OG1	1:A:243:LYS:HD2	2.19	0.42
3:C:4:TYR:CZ	4:G:102:GLY:HA2	2.55	0.42
1:D:192:HIS:O	1:D:193:PRO:C	2.63	0.42
1:D:231:VAL:HG11	1:D:244:TRP:CZ2	2.55	0.42
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.55	0.41
1:A:202:ARG:NH1	2:B:98:ASP:O	2.53	0.41
5:H:34:ARG:NH1	5:H:62:TYR:OH	2.52	0.41
5:L:132:ILE:HG23	5:L:191:ALA:HB1	2.02	0.41
1:D:103:LEU:C	1:D:110:LEU:HD12	2.45	0.41
5:H:12:VAL:HG22	5:H:114:LEU:HD21	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:TYR:CE2	3:C:9:MET:HE2	2.56	0.40
6:K:32:TYR:CE1	6:K:78:LYS:HG2	2.55	0.40
5:L:32:TRP:CD1	5:L:72:PHE:CE2	3.10	0.40
6:K:67:GLY:C	6:K:69:PHE:H	2.29	0.40

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	273/276 (99%)	258 (94%)	15 (6%)	0	100	100
1	D	273/276 (99%)	258 (94%)	15 (6%)	0	100	100
2	B	96/99 (97%)	89 (93%)	6 (6%)	1 (1%)	12	24
2	E	96/99 (97%)	90 (94%)	5 (5%)	1 (1%)	12	24
3	C	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	F	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
4	G	112/205 (55%)	100 (89%)	12 (11%)	0	100	100
5	H	224/238 (94%)	205 (92%)	19 (8%)	0	100	100
5	L	232/238 (98%)	213 (92%)	19 (8%)	0	100	100
6	K	114/160 (71%)	102 (90%)	12 (10%)	0	100	100
All	All	1434/1609 (89%)	1327 (92%)	105 (7%)	2 (0%)	48	69

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	47	PRO
2	B	47	PRO

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	234/234 (100%)	225 (96%)	9 (4%)	29	54
1	D	234/234 (100%)	225 (96%)	9 (4%)	29	54
2	B	93/94 (99%)	90 (97%)	3 (3%)	34	59
2	E	92/94 (98%)	90 (98%)	2 (2%)	45	69
3	C	7/7 (100%)	6 (86%)	1 (14%)	3	6
3	F	7/7 (100%)	7 (100%)	0	100	100
4	G	100/184 (54%)	91 (91%)	9 (9%)	9	20
5	H	197/204 (97%)	186 (94%)	11 (6%)	19	40
5	L	202/204 (99%)	193 (96%)	9 (4%)	24	48
6	K	103/144 (72%)	97 (94%)	6 (6%)	18	38
All	All	1269/1406 (90%)	1210 (95%)	59 (5%)	24	48

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	103	LEU
1	A	108	ARG
1	A	114	LEU
1	A	180	LEU
1	A	200	THR
1	A	206	LEU
1	A	223	GLU
1	A	260	ARG
1	A	266	LEU
2	B	29	GLN
2	B	75	THR
2	B	83	LYS
3	C	9	MET
1	D	103	LEU
1	D	108	ARG
1	D	114	LEU

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Mol	Chain	Res	Type
1	D	180	LEU
1	D	200	THR
1	D	206	LEU
1	D	223	GLU
1	D	260	ARG
1	D	266	LEU
2	E	4	THR
2	E	70	PHE
4	G	9	GLN
4	G	20	VAL
4	G	59	SER
4	G	71	THR
4	G	76	ARG
4	G	79	LYS
4	G	85	ILE
4	G	105	LYS
4	G	115	THR
5	H	3	THR
5	H	8	SER
5	H	52	SER
5	H	63	LYS
5	H	73	SER
5	H	99	THR
5	H	112	GLU
5	H	165	SER
5	H	180	SER
5	H	205	VAL
5	H	235	ARG
6	K	26	THR
6	K	31	SER
6	K	56	SER
6	K	70	THR
6	K	76	ARG
6	K	115	THR
5	L	52	SER
5	L	116	ASN
5	L	126	GLU
5	L	153	GLU
5	L	161	LYS
5	L	168	CYS
5	L	182	SER
5	L	206	GLN

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Mol	Chain	Res	Type
5	L	212	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	87	GLN
1	A	97	GLN
1	A	127	ASN
1	A	169	HIS
1	A	174	ASN
1	A	220	ASN
1	A	263	HIS
2	B	29	GLN
2	B	38	GLN
2	B	67	HIS
3	C	5	ASN
1	D	65	GLN
1	D	127	ASN
1	D	169	HIS
1	D	174	ASN
1	D	263	HIS
2	E	29	GLN
4	G	9	GLN
4	G	10	GLN
4	G	45	GLN
4	G	74	ASN
5	H	22	HIS
5	H	35	GLN
5	H	71	ASN
5	H	196	ASN
5	H	199	ASN
5	H	206	GLN
6	K	83	HIS
5	L	39	HIS
5	L	69	GLN
5	L	71	ASN
5	L	116	ASN
5	L	208	HIS
5	L	227	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

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	275/276 (99%)	-0.02	5 (1%) 67 60	37, 64, 119, 133	0
1	D	275/276 (99%)	0.02	3 (1%) 78 73	43, 70, 118, 142	0
2	B	98/99 (98%)	-0.32	0 100 100	39, 59, 87, 103	0
2	E	98/99 (98%)	-0.15	0 100 100	41, 66, 91, 116	0
3	C	9/9 (100%)	-0.84	0 100 100	42, 44, 57, 58	0
3	F	9/9 (100%)	-0.81	0 100 100	48, 51, 60, 63	0
4	G	114/205 (55%)	-0.27	0 100 100	41, 63, 102, 130	0
5	H	230/238 (96%)	0.12	6 (2%) 57 48	36, 80, 138, 159	0
5	L	236/238 (99%)	-0.02	3 (1%) 75 69	44, 79, 127, 151	0
6	K	118/160 (73%)	-0.15	1 (0%) 82 78	47, 70, 102, 123	0
All	All	1462/1609 (90%)	-0.06	18 (1%) 76 71	36, 70, 121, 159	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	H	193	PHE	3.5
5	H	133	ALA	3.0
5	H	221	PRO	2.9
1	A	220	ASN	2.9
1	D	107	TRP	2.9
6	K	120	ILE	2.8
5	H	215	LYS	2.8
1	A	107	TRP	2.8
5	H	190	SER	2.8
1	D	2	PRO	2.6
1	A	138	MET	2.6
5	L	217	PRO	2.3
1	D	177	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
5	L	39	HIS	2.1
5	H	137	LYS	2.1
5	L	129	LYS	2.1
1	A	173	LYS	2.0
1	A	2	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 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.