



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 29, 2022 – 12:06 AM EST

PDB ID : 7T0L
Title : HLA-B*27:05 in complex with the pan-HLA-Ia monoclonal antibody W6/32
Authors : Vivian, J.P.; Rossjohn, J.
Deposited on : 2021-11-29
Resolution : 3.00 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	:	2.31.2
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)	:	2.31.2

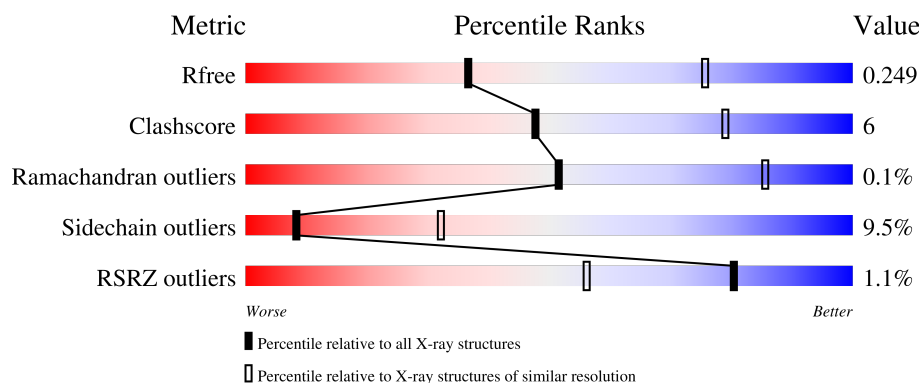
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION






The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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	 78% 20% .
1	D	276	 74% 22% .
2	B	99	 79% 20% .
2	E	99	 3% 81% 18% .
3	C	9	 89% 11%

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Mol	Chain	Length	Quality of chain
3	F	9	<div><div></div><div>78%22%</div></div>
4	G	218	<div><div>%</div><div></div><div>83%17%</div><div>.</div></div>
4	H	218	<div><div></div><div>80%17%</div><div>..</div></div>
5	K	211	<div><div>%</div><div></div><div>75%22%</div><div>.</div></div>
5	L	211	<div><div>4%</div><div></div><div>83%13%</div><div>.</div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12916 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 MHC class I antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2252	1401	409	436	6			
1	D	276	Total	C	N	O	S	0	0	0
			2252	1401	409	436	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	67	SER	CYS	engineered mutation	UNP A3F718
A	116	ASN	ASP	conflict	UNP A3F718
D	67	SER	CYS	engineered mutation	UNP A3F718
D	116	ASN	ASP	conflict	UNP A3F718

- 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	4	0
			844	541	140	160	3			
2	E	99	Total	C	N	O	S	0	4	0
			844	541	140	160	3			

- Molecule 3 is a protein called PHE-ARG-TYR-ASN-GLY-LEU-ILE-HIS-ARG peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			81	51	18	12			
3	F	9	Total	C	N	O	0	0	0
			81	51	18	12			

- Molecule 4 is a protein called IgG2a heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	G	218	Total	C	N	O	S	0	0	0
			1637	1043	268	320	6			
4	H	215	Total	C	N	O	S	0	0	0
			1613	1027	265	315	6			

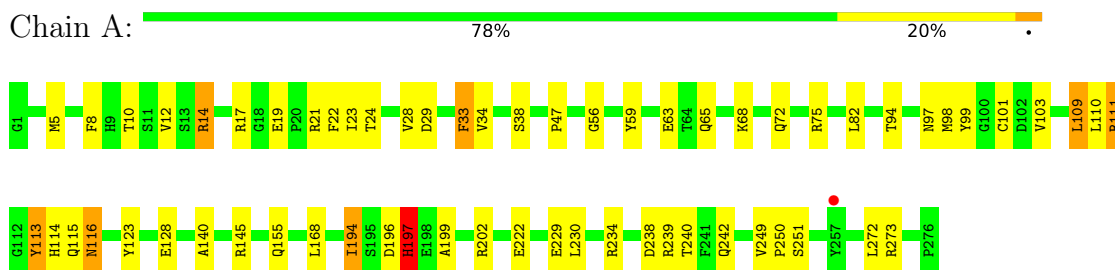
- Molecule 5 is a protein called Light chain kappa.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	K	211	Total	C	N	O	S	0	2	0
			1656	1042	273	335	6			
5	L	211	Total	C	N	O	S	0	2	0
			1656	1042	273	335	6			

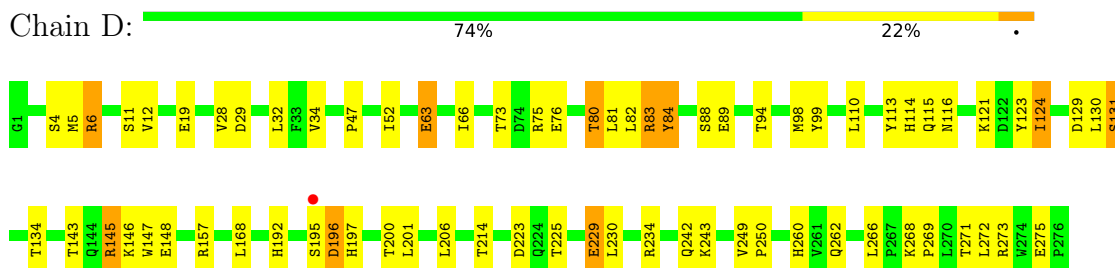
3 Residue-property plots

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.

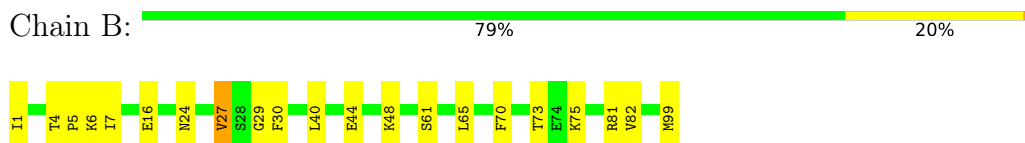
- Molecule 1: MHC class I antigen



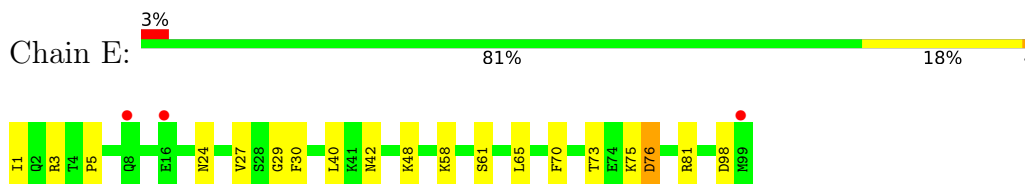
- Molecule 1: MHC class I antigen



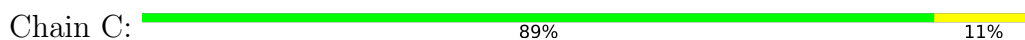
- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin

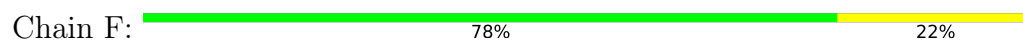


- Molecule 3: PHE-ARG-TYR-ASN-GLY-LEU-ILE-HIS-ARG peptide

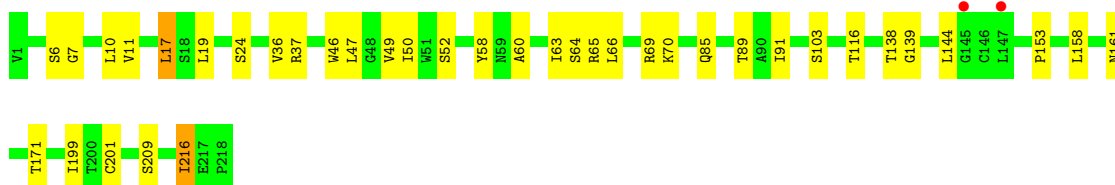
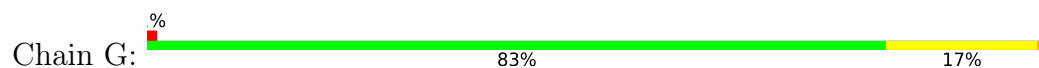




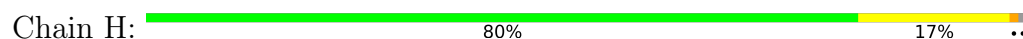
- Molecule 3: PHE-ARG-TYR-ASN-GLY-LEU-ILE-HIS-ARG peptide



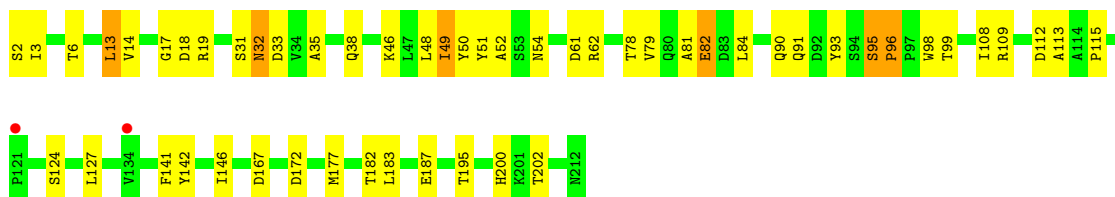
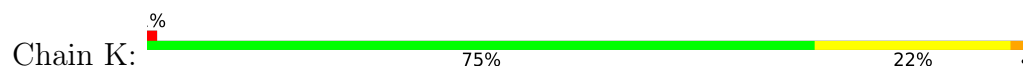
- Molecule 4: IgG2a heavy chain



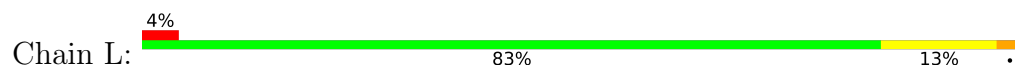
- Molecule 4: IgG2a heavy chain



- Molecule 5: Light chain kappa



- Molecule 5: Light chain kappa





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	142.78Å 146.66Å 177.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.58 – 3.00 102.31 – 3.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (42.58-3.00) 92.7 (102.31-3.00)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 3.01Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.198 , 0.252 0.213 , 0.249	Depositor DCC
R_{free} test set	1884 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	66.2	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 41.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.136 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	12916	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% 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	0.44	0/2314	0.63	1/3147 (0.0%)
1	D	0.45	0/2314	0.62	0/3147
2	B	0.41	0/879	0.63	0/1189
2	E	0.41	0/879	0.61	0/1189
3	C	0.49	0/83	0.72	0/108
3	F	0.40	0/83	0.63	0/108
4	G	0.47	0/1682	0.69	0/2304
4	H	0.45	0/1657	0.66	0/2269
5	K	0.46	0/1703	0.62	1/2319 (0.0%)
5	L	0.49	1/1703 (0.1%)	0.68	3/2319 (0.1%)
All	All	0.45	1/13297 (0.0%)	0.65	5/18099 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	L	96	PRO	N-CD	5.65	1.55	1.47

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	L	96	PRO	C-N-CD	6.75	142.57	128.40
5	K	96	PRO	C-N-CD	6.42	141.89	128.40
1	A	197	HIS	N-CA-C	6.00	127.20	111.00
5	L	95	SER	C-N-CD	5.07	139.05	128.40
5	L	69	GLY	N-CA-C	5.04	125.71	113.10

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	2252	0	2111	32	0
1	D	2252	0	2111	38	0
2	B	844	0	824	10	0
2	E	844	0	824	6	0
3	C	81	0	75	1	0
3	F	81	0	75	2	0
4	G	1637	0	1606	15	0
4	H	1613	0	1582	16	0
5	K	1656	0	1582	27	0
5	L	1656	0	1582	18	0
All	All	12916	0	12372	153	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 (153) 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:197:HIS:O	1:A:250:PRO:HA	1.71	0.91
1:D:81:LEU:HA	1:D:84:TYR:HB2	1.53	0.87
5:L:14:VAL:HG21	5:L:20:VAL:HG11	1.57	0.84
1:A:197:HIS:O	1:A:251:SER:N	2.17	0.78
5:K:3:ILE:O	5:K:99:THR:HG21	1.85	0.77
1:A:202:ARG:HH12	2:B:99:MET:HG2	1.49	0.77
4:G:158:LEU:HD11	4:G:201:CYS:HB2	1.67	0.76
4:H:158:LEU:HD11	4:H:201:CYS:HB2	1.68	0.75
1:D:234:ARG:HG3	1:D:242:GLN:HG3	1.70	0.73
1:D:230:LEU:HD22	1:D:243:LYS:HE3	1.71	0.72
4:H:10:LEU:HB2	4:H:153:PRO:HG3	1.75	0.69
5:K:38:GLN:HB2	5:K:48:LEU:HD11	1.74	0.69
4:H:89:THR:HG23	4:H:116:THR:HA	1.74	0.69
1:A:238:ASP:O	1:A:239:ARG:HB2	1.96	0.65
5:L:183:LEU:HD22	5:L:187:GLU:HG2	1.80	0.64
5:L:38:GLN:HB2	5:L:48:LEU:HD11	1.81	0.62
1:D:5:MET:HB2	1:D:168:LEU:HD13	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:214:THR:HB	1:D:262:GLN:HB2	1.82	0.61
2:E:24:ASN:HB3	2:E:65:LEU:HD11	1.83	0.60
1:D:201:LEU:HD21	1:D:273:ARG:HH22	1.65	0.60
1:A:111:ARG:HE	1:A:128:GLU:HG2	1.67	0.60
4:H:104:ALA:HB3	5:L:92:ASP:HB3	1.83	0.60
5:L:119:ILE:HD12	5:L:196:CYS:HB2	1.83	0.60
1:D:99:TYR:HB3	1:D:114:HIS:CD2	2.36	0.60
5:K:17:GLY:HA2	5:K:78:THR:HG23	1.83	0.60
1:D:83:ARG:HG2	1:D:83:ARG:NH1	2.17	0.59
1:A:197:HIS:O	1:A:250:PRO:CA	2.49	0.59
1:D:145:ARG:O	1:D:147:TRP:N	2.37	0.58
4:G:7:GLY:HA3	4:G:19:LEU:HD23	1.86	0.58
1:D:83:ARG:CG	1:D:83:ARG:HH11	2.16	0.57
1:A:202:ARG:NH1	2:B:99:MET:HG2	2.19	0.57
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.87	0.55
1:A:19:GLU:HB3	1:A:75:ARG:HE	1.70	0.55
5:K:51:TYR:O	5:K:52:ALA:HB3	2.07	0.55
5:K:124:SER:HA	5:K:127:LEU:HD12	1.89	0.55
1:D:260:HIS:HD1	1:D:271:THR:HG22	1.70	0.55
5:K:50:TYR:O	5:K:51:TYR:HB2	2.05	0.55
5:K:32:ASN:O	5:K:32:ASN:ND2	2.39	0.55
1:A:65:GLN:HB2	4:G:85:GLN:OE1	2.06	0.55
1:D:63:GLU:HA	1:D:66:ILE:HD12	1.88	0.55
1:D:196:ASP:HB3	1:D:197:HIS:ND1	2.23	0.54
4:G:58:TYR:HB2	4:G:63:ILE:HD13	1.89	0.54
5:K:2:SER:O	5:K:3:ILE:C	2.46	0.53
1:D:145:ARG:O	1:D:146:LYS:C	2.47	0.53
4:G:89:THR:HG23	4:G:116:THR:HA	1.89	0.53
4:G:11:VAL:HG21	4:G:17:LEU:HB2	1.91	0.53
5:L:13:LEU:HD12	5:L:107:GLU:HB3	1.92	0.52
5:K:96:PRO:O	5:K:98:TRP:CD2	2.63	0.52
5:L:135:VAL:HG22	5:L:180:THR:HG23	1.91	0.51
4:G:36:VAL:HG22	4:G:46:TRP:HA	1.93	0.51
4:G:37:ARG:HG2	4:G:47:LEU:HD21	1.93	0.51
5:L:30:VAL:HG12	5:L:93:TYR:HB3	1.92	0.51
1:A:229:GLU:HG2	4:H:69:ARG:HH21	1.75	0.51
5:L:14:VAL:HG21	5:L:20:VAL:CG1	2.35	0.50
4:H:58:TYR:HB2	4:H:63:ILE:HD13	1.93	0.50
1:A:111:ARG:HD2	1:A:113:TYR:CD2	2.46	0.50
5:K:18:ASP:O	5:K:79:VAL:HG23	2.11	0.50
1:D:12:VAL:HG22	1:D:94:THR:HG23	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:L:29:SER:HA	5:L:69:GLY:O	2.11	0.50
1:D:116:ASN:OD1	1:D:124:ILE:HG22	2.12	0.50
5:K:51:TYR:HB2	5:K:54:ASN:HD22	1.75	0.50
2:B:4:THR:HG22	5:L:96:PRO:HD3	1.94	0.50
2:B:7:ILE:HG12	2:B:82:VAL:HG21	1.94	0.50
5:K:13:LEU:HD23	5:K:109:ARG:HH21	1.77	0.49
1:A:197:HIS:N	1:A:197:HIS:ND1	2.60	0.49
5:K:81:ALA:HA	5:K:108:ILE:HD13	1.94	0.49
5:L:51:TYR:O	5:L:52:ALA:HB3	2.12	0.49
5:K:31:SER:O	5:K:31:SER:OG	2.26	0.49
1:A:28:VAL:O	1:A:29:ASP:HB2	2.12	0.49
1:D:123:TYR:CE2	1:D:143:THR:HG21	2.48	0.49
1:D:229:GLU:HG2	4:G:69:ARG:HH21	1.78	0.49
1:D:83:ARG:HG2	1:D:83:ARG:HH11	1.77	0.49
4:G:144:LEU:HD22	4:G:216:ILE:HG21	1.95	0.49
5:K:115:PRO:HB3	5:K:141:PHE:CD1	2.47	0.49
1:A:155:GLN:HG2	3:C:3:TYR:HE1	1.77	0.48
4:H:125:PRO:HB3	4:H:151:TYR:HB3	1.94	0.48
1:A:14:ARG:HD3	1:A:17:ARG:HB3	1.94	0.48
4:G:10:LEU:HB2	4:G:153:PRO:HG3	1.95	0.48
4:H:161:ASN:HD22	4:H:165:LEU:HD11	1.78	0.48
4:H:60:ALA:O	4:H:63:ILE:HG12	2.13	0.48
5:K:82:GLU:H	5:K:82:GLU:HG3	1.43	0.48
1:A:194:ILE:HG13	1:A:199:ALA:HA	1.95	0.48
5:K:95:SER:HB3	5:K:96:PRO:HD2	1.94	0.48
5:L:91:GLN:O	5:L:98:TRP:HB3	2.14	0.48
1:A:12:VAL:HG12	1:A:94:THR:HG23	1.96	0.48
2:B:40:LEU:HD11	2:B:81:ARG:HB2	1.96	0.48
1:D:192:HIS:HB2	1:D:200:THR:HB	1.94	0.48
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.96	0.48
1:D:83:ARG:NH1	1:D:83:ARG:CG	2.73	0.48
1:D:268:LYS:HG2	1:D:269:PRO:HD2	1.96	0.48
5:K:146:ILE:HD12	5:K:200:HIS:CD2	2.49	0.47
5:L:95:SER:HB3	5:L:96:PRO:CD	2.43	0.47
1:D:66:ILE:HD13	3:F:2:ARG:HB2	1.95	0.47
5:K:183:LEU:HD22	5:K:187:GLU:HG2	1.95	0.47
1:D:81:LEU:CA	1:D:84:TYR:HB2	2.36	0.47
1:D:145:ARG:O	1:D:148:GLU:N	2.44	0.47
1:D:196:ASP:HB3	1:D:197:HIS:H	1.47	0.47
1:D:6:ARG:HB3	1:D:98:MET:HE3	1.96	0.47
1:D:195:SER:O	1:D:196:ASP:C	2.52	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:ARG:HG3	1:A:242:GLN:HG3	1.97	0.47
1:D:80:THR:O	1:D:84:TYR:N	2.45	0.46
1:D:99:TYR:HB3	1:D:114:HIS:HD2	1.77	0.46
2:E:40:LEU:HD11	2:E:81:ARG:HB2	1.97	0.46
1:A:33:PHE:H	1:A:33:PHE:HD1	1.63	0.46
2:B:5:PRO:HB2	2:B:27:VAL:HG12	1.97	0.46
2:B:29:GLY:HA2	2:B:61[B]:SER:HB3	1.98	0.46
1:D:47:PRO:HB3	1:D:52:ILE:HG23	1.97	0.46
5:K:95:SER:O	5:K:96:PRO:C	2.54	0.46
1:A:123:TYR:CZ	1:A:140:ALA:HA	2.51	0.45
2:E:5:PRO:HB3	2:E:30:PHE:HB3	1.97	0.45
1:A:101:CYS:HB2	1:A:109:LEU:HD12	1.98	0.45
4:H:148:VAL:HB	4:H:183:LEU:HB3	1.98	0.45
4:G:50:ILE:HG12	4:G:70:LYS:HE2	1.99	0.45
5:K:113:ALA:C	5:K:202:THR:HG21	2.37	0.45
2:B:4:THR:HG22	5:L:96:PRO:CD	2.47	0.44
1:D:129:ASP:O	1:D:131:SER:OG	2.33	0.44
2:E:29:GLY:HA2	2:E:61[B]:SER:HB3	1.99	0.44
4:H:129:PRO:HD3	4:H:214:LYS:HE3	2.00	0.44
1:D:201:LEU:HD22	1:D:249:VAL:HG11	1.99	0.44
5:L:3:ILE:HG23	5:L:28:GLN:H	1.81	0.44
4:H:130:LEU:HD11	4:H:147:LEU:HB2	2.00	0.43
5:L:90:GLN:HG2	5:L:91:GLN:N	2.33	0.43
1:A:145:ARG:NH2	2:E:98:ASP:OD2	2.48	0.43
5:L:12:LEU:HD21	5:L:20:VAL:HB	2.00	0.43
1:A:97:ASN:HD22	1:A:116:ASN:HB3	1.84	0.43
2:B:24:ASN:HB3	2:B:65:LEU:HD11	2.00	0.42
1:A:8:PHE:O	1:A:24:THR:HA	2.20	0.42
5:K:32:ASN:ND2	5:K:32:ASN:C	2.73	0.42
1:D:197:HIS:O	1:D:250:PRO:HA	2.19	0.42
4:H:66:LEU:HD23	4:H:67:SER:N	2.35	0.42
5:K:95:SER:CB	5:K:96:PRO:HD2	2.49	0.42
1:A:99:TYR:HB3	1:A:114:HIS:ND1	2.35	0.42
1:A:197:HIS:O	1:A:250:PRO:C	2.58	0.41
1:A:21:ARG:HE	1:A:23:ILE:HD11	1.85	0.41
4:G:60:ALA:HA	4:G:63:ILE:CG1	2.51	0.41
1:A:10:THR:O	1:A:22:PHE:HA	2.21	0.41
4:H:17:LEU:O	4:H:80:LYS:HA	2.20	0.41
1:D:145:ARG:C	1:D:147:TRP:N	2.71	0.41
5:K:112:ASP:HA	5:K:142:TYR:HB3	2.03	0.41
1:D:73:THR:HG23	3:F:8:HIS:CE1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:ASP:HB3	1:A:197:HIS:H	1.49	0.41
1:D:28:VAL:O	1:D:29:ASP:HB2	2.21	0.41
4:G:138:THR:HB	4:G:139:GLY:H	1.55	0.41
1:D:19:GLU:HB3	1:D:75:ARG:HE	1.85	0.41
4:H:60:ALA:C	4:H:62:PHE:H	2.24	0.41
5:K:49:ILE:HA	5:K:54:ASN:O	2.20	0.41
1:A:56:GLY:O	1:A:59:TYR:HB3	2.21	0.40
2:E:42:ASN:HD21	2:E:76:ASP:HA	1.85	0.40
5:K:33:ASP:HB2	5:K:93:TYR:HB2	2.03	0.40
1:A:22:PHE:HB3	1:A:38:SER:HB3	2.03	0.40
4:G:46:TRP:HZ2	4:G:49:VAL:HG12	1.86	0.40
4:H:156:VAL:HG12	4:H:205:HIS:HB2	2.02	0.40
5:K:35:ALA:HB3	5:K:90:GLN:HB3	2.04	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	274/276 (99%)	256 (93%)	17 (6%)	1 (0%)	34	72
1	D	274/276 (99%)	254 (93%)	20 (7%)	0	100	100
2	B	101/99 (102%)	98 (97%)	3 (3%)	0	100	100
2	E	101/99 (102%)	98 (97%)	3 (3%)	0	100	100
3	C	7/9 (78%)	5 (71%)	2 (29%)	0	100	100
3	F	7/9 (78%)	7 (100%)	0	0	100	100
4	G	216/218 (99%)	207 (96%)	9 (4%)	0	100	100
4	H	213/218 (98%)	197 (92%)	15 (7%)	1 (0%)	29	68
5	K	211/211 (100%)	203 (96%)	8 (4%)	0	100	100
5	L	211/211 (100%)	202 (96%)	9 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1615/1626 (99%)	1527 (95%)	86 (5%)	2 (0%)	51 85

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	H	61	ALA
1	A	194	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	235/235 (100%)	212 (90%)	23 (10%)	8 30
1	D	235/235 (100%)	204 (87%)	31 (13%)	4 18
2	B	98/94 (104%)	89 (91%)	9 (9%)	9 34
2	E	98/94 (104%)	89 (91%)	9 (9%)	9 34
3	C	7/8 (88%)	7 (100%)	0	100 100
3	F	7/8 (88%)	7 (100%)	0	100 100
4	G	187/187 (100%)	173 (92%)	14 (8%)	13 43
4	H	184/187 (98%)	169 (92%)	15 (8%)	11 39
5	K	189/187 (101%)	171 (90%)	18 (10%)	8 32
5	L	189/187 (101%)	174 (92%)	15 (8%)	12 41
All	All	1429/1422 (100%)	1295 (91%)	134 (9%)	8 32

All (134) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	33	PHE
1	A	34	VAL
1	A	47	PRO
1	A	63	GLU

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Mol	Chain	Res	Type
1	A	68	LYS
1	A	72	GLN
1	A	82	LEU
1	A	98	MET
1	A	103	VAL
1	A	109	LEU
1	A	110	LEU
1	A	111	ARG
1	A	113	TYR
1	A	115	GLN
1	A	116	ASN
1	A	197	HIS
1	A	222	GLU
1	A	230	LEU
1	A	240	THR
1	A	249	VAL
1	A	272	LEU
1	A	273	ARG
2	B	1	ILE
2	B	6	LYS
2	B	16	GLU
2	B	27	VAL
2	B	44	GLU
2	B	48	LYS
2	B	70	PHE
2	B	73	THR
2	B	75	LYS
1	D	4	SER
1	D	6	ARG
1	D	11	SER
1	D	32	LEU
1	D	34	VAL
1	D	63	GLU
1	D	76	GLU
1	D	80	THR
1	D	82	LEU
1	D	83	ARG
1	D	84	TYR
1	D	88	SER
1	D	89	GLU
1	D	110	LEU
1	D	113	TYR

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Mol	Chain	Res	Type
1	D	115	GLN
1	D	121	LYS
1	D	124	ILE
1	D	130	LEU
1	D	131	SER
1	D	134	THR
1	D	145	ARG
1	D	157	ARG
1	D	196	ASP
1	D	206	LEU
1	D	223	ASP
1	D	225	THR
1	D	229	GLU
1	D	266	LEU
1	D	272	LEU
1	D	275	GLU
2	E	1	ILE
2	E	3	ARG
2	E	27	VAL
2	E	48	LYS
2	E	58	LYS
2	E	70	PHE
2	E	73	THR
2	E	75	LYS
2	E	76	ASP
4	G	6	SER
4	G	17	LEU
4	G	24	SER
4	G	52	SER
4	G	64	SER
4	G	65	ARG
4	G	66	LEU
4	G	91	ILE
4	G	103	SER
4	G	161	ASN
4	G	171	THR
4	G	199	ILE
4	G	209	SER
4	G	216	ILE
4	H	24	SER
4	H	45	GLU
4	H	65	ARG

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Mol	Chain	Res	Type
4	H	66	LEU
4	H	85	GLN
4	H	91	ILE
4	H	100	THR
4	H	115	VAL
4	H	154	GLU
4	H	165	LEU
4	H	166	SER
4	H	183	LEU
4	H	202	ASN
4	H	213	ASP
4	H	215	LYS
5	K	6	THR
5	K	13	LEU
5	K	14	VAL
5	K	19	ARG
5	K	32	ASN
5	K	46	LYS
5	K	49	ILE
5	K	61	ASP
5	K	62	ARG
5	K	82	GLU
5	K	84	LEU
5	K	91	GLN
5	K	95	SER
5	K	167	ASP
5	K	172	ASP
5	K	177	MET
5	K	182	THR
5	K	195	THR
5	L	5	MET
5	L	13	LEU
5	L	14	VAL
5	L	19	ARG
5	L	21	THR
5	L	44	SER
5	L	49	ILE
5	L	64	THR
5	L	84	LEU
5	L	91	GLN
5	L	92	ASP
5	L	96	PRO

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Mol	Chain	Res	Type
5	L	112	ASP
5	L	177	MET
5	L	180	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	96	GLN
1	D	115	GLN
1	D	144	GLN
4	H	85	GLN
5	K	54	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	276/276 (100%)	-0.53	1 (0%) 92 79	52, 84, 124, 142	0
1	D	276/276 (100%)	-0.49	1 (0%) 92 79	48, 99, 135, 151	0
2	B	99/99 (100%)	-0.44	0 100 100	58, 83, 139, 149	0
2	E	99/99 (100%)	-0.16	3 (3%) 50 22	60, 91, 149, 159	0
3	C	9/9 (100%)	-0.52	0 100 100	45, 66, 69, 73	0
3	F	9/9 (100%)	-0.13	0 100 100	84, 91, 95, 98	0
4	G	218/218 (100%)	-0.40	2 (0%) 84 63	31, 59, 153, 193	0
4	H	215/218 (98%)	-0.36	1 (0%) 91 75	41, 67, 144, 163	0
5	K	211/211 (100%)	-0.35	2 (0%) 84 63	48, 106, 166, 184	0
5	L	211/211 (100%)	-0.28	8 (3%) 40 16	55, 103, 166, 184	0
All	All	1623/1626 (99%)	-0.40	18 (1%) 80 56	31, 89, 149, 193	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	G	145	GLY	4.7
5	K	121	PRO	3.1
5	L	137	PHE	2.8
2	E	99	MET	2.5
5	K	134	VAL	2.3
5	L	120	PHE	2.3
5	L	2	SER	2.3
1	D	195	SER	2.2
1	A	257	TYR	2.2
2	E	8	GLN	2.2
4	G	147	LEU	2.2
5	L	71	ASP	2.1
5	L	32	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
4	H	144	LEU	2.1
5	L	163	ASN	2.1
5	L	211	PHE	2.1
2	E	16	GLU	2.1
5	L	119	ILE	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 [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.