



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:20 AM UTC

PDB ID : 3OX8 / pdb_00003ox8
Title : Crystal Structure of HLA A*02:03 Bound to HBV Core 18-27
Authors : Liu, J.; Chen, Y.; Lai, L.; Ren, E.
Deposited on : 2010-09-21
Resolution : 2.16 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

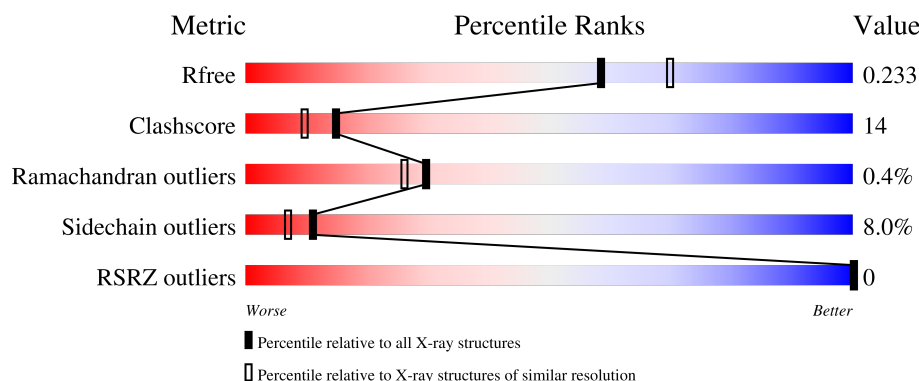
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

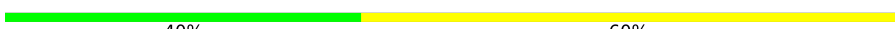
The reported resolution of this entry is 2.16 Å.

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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	275	 73% 23% .
1	D	275	 73% 24% .
2	B	100	 82% 16% .
2	E	100	 79% 15% 6%
3	C	10	 40% 60%

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Mol	Chain	Length	Quality of chain
3	F	10	 70% 30%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6820 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	275	Total	C	N	O	S	0	1	0
			2265	1414	413	429	9			
1	D	275	Total	C	N	O	S	0	1	0
			2265	1414	413	429	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 10mer peptide from Pre-core-protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	10	Total	C	N	O	0	0	0
			83	58	10	15			
3	F	10	Total	C	N	O	0	0	0
			83	58	10	15			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	165	Total	O	0	0
			165	165		

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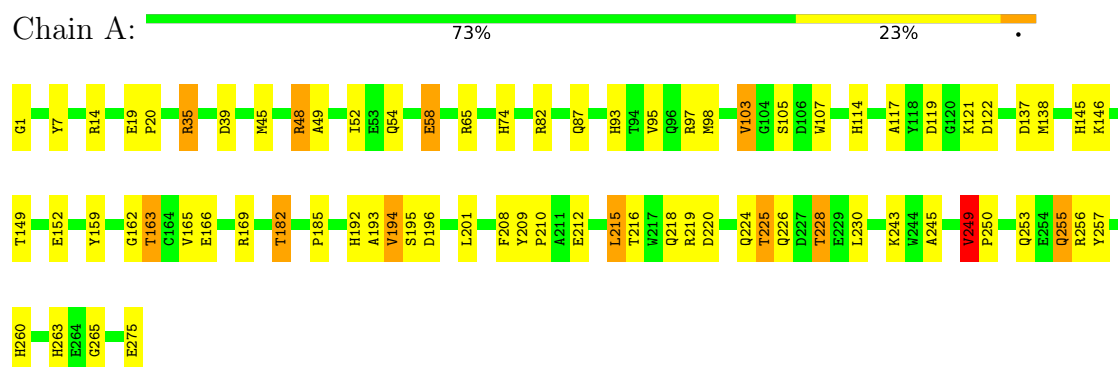
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	60	Total 60	O 60	0	0
4	C	4	Total 4	O 4	0	0
4	D	161	Total 161	O 161	0	0
4	E	55	Total 55	O 55	0	0
4	F	5	Total 5	O 5	0	0

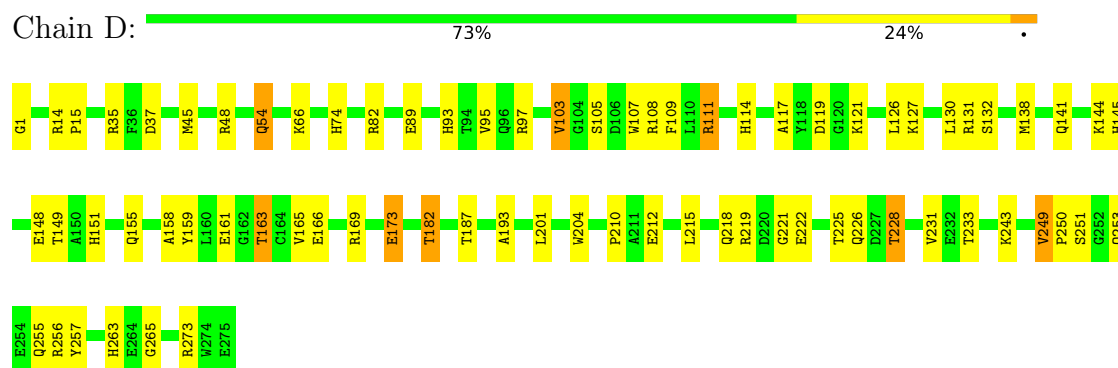
3 Residue-property plots

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.

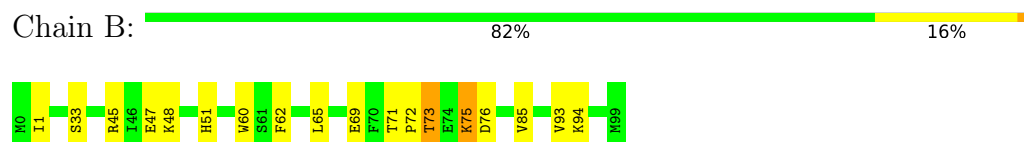
• Molecule 1: MHC class I antigen



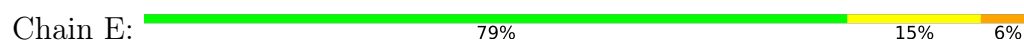
• Molecule 1: MHC class I antigen



• Molecule 2: Beta-2-microglobulin



• Molecule 2: Beta-2-microglobulin

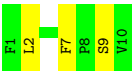




- Molecule 3: 10mer peptide from Pre-core-protein



- Molecule 3: 10mer peptide from Pre-core-protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.28Å 68.27Å 68.32Å 70.20° 84.40° 84.45°	Depositor
Resolution (Å)	64.12 – 2.16 64.12 – 2.16	Depositor EDS
% Data completeness (in resolution range)	95.9 (64.12-2.16) 97.6 (64.12-2.16)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.182 , 0.234 0.182 , 0.233	Depositor DCC
R_{free} test set	2902 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	17.9	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 18.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.479 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6820	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.88% 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	0.74	0/2335	0.87	2/3168 (0.1%)
1	D	0.71	0/2335	0.84	0/3168
2	B	0.68	0/860	0.87	0/1162
2	E	0.67	0/860	0.92	0/1162
3	C	0.70	0/87	0.79	0/117
3	F	0.74	0/87	0.81	0/117
All	All	0.71	0/6564	0.86	2/8894 (0.0%)

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	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	249	VAL	CA-C-N	-5.19	114.57	119.76
1	A	249	VAL	C-N-CA	-5.19	114.57	119.76

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	111	ARG	Sidechain

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	2265	0	2107	74	2
1	D	2265	0	2107	69	3
2	B	837	0	803	11	0
2	E	837	0	803	24	0
3	C	83	0	77	5	0
3	F	83	0	77	4	0
4	A	165	0	0	12	1
4	B	60	0	0	1	0
4	C	4	0	0	0	0
4	D	161	0	0	5	0
4	E	55	0	0	2	0
4	F	5	0	0	0	0
All	All	6820	0	5974	170	3

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 14.

All (170) 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:182:THR:HG21	1:D:265:GLY:HA2	1.17	1.16
1:D:182:THR:HG21	1:D:265:GLY:CA	1.83	1.08
1:D:253:GLN:NE2	1:D:256:ARG:HE	1.66	0.94
1:D:253:GLN:HE21	1:D:256:ARG:HE	1.00	0.93
1:D:54:GLN:HE21	1:D:54:GLN:HA	1.36	0.89
1:A:93:HIS:HD2	1:A:119:ASP:OD2	1.55	0.89
1:D:249:VAL:HG22	1:D:257:TYR:CE1	2.08	0.88
1:D:182:THR:CG2	1:D:265:GLY:CA	2.52	0.87
1:A:253:GLN:HE21	1:A:256:ARG:HE	1.22	0.83
1:A:35:ARG:CZ	4:A:327:HOH:O	2.27	0.83
1:A:249:VAL:HG22	1:A:257:TYR:CE1	2.15	0.82
1:A:145:HIS:O	1:A:149:THR:HG23	1.80	0.81
2:E:73:THR:CG2	2:E:76:ASP:H	1.93	0.80
2:B:73:THR:CG2	2:B:76:ASP:H	1.94	0.79
1:A:226:GLN:HG2	4:A:294:HOH:O	1.82	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:73:THR:HG22	2:B:76:ASP:H	1.48	0.79
1:D:127:LYS:HE2	1:D:132:SER:OG	1.83	0.79
1:D:97:ARG:HH11	1:D:114:HIS:HE1	1.29	0.78
1:D:93:HIS:HD2	1:D:119:ASP:OD2	1.66	0.78
1:D:253:GLN:HE21	1:D:256:ARG:NE	1.81	0.78
1:A:255:GLN:HG2	4:A:300:HOH:O	1.84	0.77
1:A:263:HIS:CD2	1:A:265:GLY:H	2.02	0.77
1:A:97:ARG:HD3	1:A:114:HIS:HE1	1.49	0.76
2:E:73:THR:HG22	2:E:76:ASP:H	1.51	0.76
1:D:263:HIS:HD2	1:D:265:GLY:H	1.32	0.75
1:A:253:GLN:NE2	1:A:256:ARG:HE	1.86	0.74
1:D:182:THR:CG2	1:D:265:GLY:HA3	2.17	0.74
1:A:263:HIS:HD2	1:A:265:GLY:H	1.34	0.72
1:D:249:VAL:HG22	1:D:257:TYR:CZ	2.23	0.72
2:E:74:GLU:N	2:E:75:LYS:HZ2	1.88	0.72
1:A:196:ASP:HB2	4:A:353:HOH:O	1.89	0.71
1:A:97:ARG:HH11	1:A:114:HIS:CE1	2.07	0.71
1:A:35:ARG:HH11	1:A:48:ARG:NH1	1.89	0.71
2:B:75:LYS:H	2:B:75:LYS:HZ2	1.35	0.70
1:A:103:VAL:HG13	1:A:107:TRP:HA	1.73	0.70
1:D:263:HIS:CD2	1:D:265:GLY:H	2.07	0.70
2:E:17:ASN:HD21	2:E:74:GLU:HG3	1.56	0.70
1:D:253:GLN:NE2	1:D:256:ARG:NE	2.40	0.69
2:E:75:LYS:N	2:E:75:LYS:HE3	2.08	0.68
1:A:14:ARG:HD3	4:A:355:HOH:O	1.91	0.68
1:D:182:THR:CG2	1:D:265:GLY:HA2	2.07	0.68
1:D:97:ARG:HH11	1:D:114:HIS:CE1	2.11	0.68
1:A:82:ARG:NH1	1:A:87:GLN:O	2.27	0.67
1:D:141:GLN:OE1	1:D:145:HIS:CE1	2.48	0.67
1:D:54:GLN:HE21	1:D:54:GLN:CA	2.07	0.66
1:D:117:ALA:HB2	2:E:60:TRP:CE2	2.31	0.66
1:A:97:ARG:HH11	1:A:114:HIS:HE1	1.43	0.65
2:B:73:THR:HG21	2:B:76:ASP:OD2	1.97	0.65
1:A:182:THR:HG21	1:A:265:GLY:HA2	1.79	0.65
1:A:93:HIS:CD2	1:A:119:ASP:OD2	2.45	0.64
1:A:225:THR:O	1:A:228:THR:HB	1.98	0.64
1:D:145:HIS:O	1:D:149:THR:HG23	1.98	0.64
1:A:97:ARG:HD3	1:A:114:HIS:CE1	2.33	0.63
1:D:225:THR:O	1:D:228:THR:HB	1.99	0.63
2:E:74:GLU:H	2:E:75:LYS:HZ2	1.46	0.62
1:A:249:VAL:HG22	1:A:257:TYR:CZ	2.33	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:73:THR:HG21	2:E:76:ASP:CG	2.24	0.62
1:D:273:ARG:NH1	4:D:711:HOH:O	2.29	0.60
1:D:231:VAL:O	1:D:243:LYS:HE2	2.01	0.60
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.37	0.59
1:A:165:VAL:O	1:A:169:ARG:HG3	2.01	0.59
2:E:73:THR:HG21	2:E:76:ASP:OD2	2.03	0.59
1:A:169:ARG:NH1	1:D:166:GLU:OE2	2.36	0.58
1:D:144:LYS:O	1:D:148:GLU:HG3	2.03	0.58
2:E:75:LYS:HE3	2:E:75:LYS:H	1.66	0.58
1:D:210:PRO:O	1:D:263:HIS:HE1	1.87	0.58
1:A:219:ARG:HH11	1:A:256:ARG:HD2	1.69	0.58
1:A:182:THR:CG2	1:A:265:GLY:HA2	2.34	0.57
1:D:35:ARG:NE	4:D:392:HOH:O	2.38	0.56
1:A:218:GLN:HG3	1:A:260:HIS:CD2	2.40	0.56
1:A:253:GLN:NE2	1:A:256:ARG:HH21	2.03	0.56
1:D:35:ARG:NH1	1:D:48:ARG:CZ	2.68	0.56
1:A:137:ASP:C	1:A:138:MET:HE2	2.29	0.56
2:B:94:LYS:NZ	4:B:635:HOH:O	2.39	0.56
2:E:22:PHE:CE2	2:E:69:GLU:HG2	2.41	0.56
2:B:73:THR:HG21	2:B:76:ASP:CG	2.31	0.56
1:A:210:PRO:O	1:A:263:HIS:HE1	1.91	0.54
1:A:35:ARG:NH1	1:A:48:ARG:NH1	2.56	0.54
1:D:114:HIS:C	1:D:114:HIS:CD2	2.85	0.54
1:D:218:GLN:HE21	1:D:221:GLY:HA2	1.72	0.54
1:A:35:ARG:NH1	1:A:48:ARG:HH12	2.06	0.54
1:A:253:GLN:HE22	1:A:256:ARG:HH21	1.56	0.54
1:A:65[A]:ARG:HD3	1:D:151:HIS:CE1	2.42	0.53
1:D:93:HIS:HE1	4:D:326:HOH:O	1.92	0.53
1:D:35:ARG:HH12	1:D:48:ARG:CZ	2.23	0.52
1:D:121:LYS:HG2	2:E:1:ILE:HD13	1.91	0.52
2:E:34:ASP:HB2	4:E:167:HOH:O	2.09	0.52
1:A:14:ARG:CD	4:A:355:HOH:O	2.53	0.52
1:A:14:ARG:NH2	1:A:39:ASP:OD2	2.41	0.52
2:E:85:VAL:HG22	4:E:106:HOH:O	2.09	0.52
1:A:138:MET:HE2	1:A:138:MET:N	2.26	0.51
1:A:165:VAL:HG11	1:D:166:GLU:HG3	1.92	0.50
1:A:169:ARG:NH1	1:D:169:ARG:NH1	2.59	0.50
1:D:249:VAL:CG2	1:D:257:TYR:CE1	2.91	0.50
1:A:220:ASP:OD1	1:A:256:ARG:HB2	2.11	0.50
2:E:73:THR:HG23	2:E:75:LYS:H	1.78	0.49
1:D:256:ARG:NH2	4:D:358:HOH:O	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:17:ASN:ND2	2:E:74:GLU:HG3	2.27	0.49
1:D:54:GLN:HA	1:D:54:GLN:NE2	2.17	0.48
1:D:187:THR:HA	1:D:204:TRP:O	2.13	0.48
1:A:14:ARG:HH22	1:A:39:ASP:CG	2.22	0.48
1:D:93:HIS:CD2	1:D:119:ASP:OD2	2.57	0.48
1:D:103:VAL:HG22	1:D:107:TRP:C	2.39	0.47
1:A:58:GLU:HB3	4:A:721:HOH:O	2.13	0.47
1:A:219:ARG:NH1	1:A:256:ARG:HD2	2.29	0.47
1:A:185:PRO:HA	1:A:208:PHE:HB3	1.98	0.46
1:D:54:GLN:CA	1:D:54:GLN:NE2	2.75	0.46
1:A:93:HIS:HE1	4:A:302:HOH:O	1.99	0.46
3:C:4:SER:OG	1:D:155:GLN:HG2	2.16	0.46
1:A:152:GLU:HG3	3:C:8:PRO:HG3	1.98	0.46
1:A:192:HIS:O	1:A:194:VAL:N	2.48	0.45
3:C:7:PHE:N	3:C:7:PHE:CD2	2.83	0.45
1:D:218:GLN:HG2	1:D:222:GLU:C	2.41	0.45
1:A:162:GLY:O	1:A:165:VAL:HG12	2.16	0.45
1:D:74:HIS:HD2	1:D:95:VAL:HG11	1.82	0.45
1:A:169:ARG:HH12	1:D:169:ARG:NH1	2.15	0.45
3:F:7:PHE:N	3:F:7:PHE:CD2	2.84	0.45
2:B:33:SER:HB3	2:B:62:PHE:CE2	2.52	0.45
1:D:126:LEU:HG	1:D:130:LEU:HA	1.98	0.45
2:E:74:GLU:H	2:E:75:LYS:NZ	2.13	0.45
1:A:121:LYS:O	1:A:122:ASP:C	2.58	0.44
3:C:6:PHE:C	3:C:6:PHE:CD1	2.95	0.44
1:A:215:LEU:HD22	1:A:243:LYS:HD3	2.00	0.44
1:A:45:MET:CE	3:C:2:LEU:HD11	2.48	0.44
1:A:255:GLN:CG	4:A:300:HOH:O	2.51	0.44
2:B:71:THR:HA	2:B:72:PRO:HD2	1.90	0.44
1:D:37:ASP:OD1	1:D:37:ASP:C	2.61	0.44
1:D:233:THR:OG1	1:D:243:LYS:HE3	2.17	0.44
1:A:159:TYR:HA	1:A:163:THR:HG23	1.99	0.44
1:D:35:ARG:CD	4:D:392:HOH:O	2.65	0.44
1:A:97:ARG:CD	1:A:114:HIS:HE1	2.24	0.43
1:D:45:MET:CE	3:F:2:LEU:HD11	2.48	0.43
1:D:97:ARG:NH1	1:D:114:HIS:HE1	2.06	0.43
1:A:209:TYR:CD1	1:A:209:TYR:C	2.96	0.43
1:D:1:GLY:O	1:D:105:SER:HA	2.18	0.43
1:D:66:LYS:HE3	3:F:2:LEU:HB2	2.00	0.43
1:D:159:TYR:HA	1:D:163:THR:HG23	2.01	0.43
1:A:58:GLU:OE2	1:D:131:ARG:HD3	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:THR:HG21	1:D:158:ALA:HB1	2.00	0.43
2:E:73:THR:HG23	2:E:75:LYS:HZ2	1.84	0.43
1:A:19:GLU:HB2	1:A:20:PRO:HD2	2.01	0.43
1:D:14:ARG:HA	1:D:15:PRO:HD2	1.89	0.43
1:A:1:GLY:O	1:A:105:SER:HA	2.18	0.43
1:A:166:GLU:OE2	4:A:343:HOH:O	2.20	0.43
1:D:103:VAL:HG22	1:D:108:ARG:C	2.43	0.43
1:D:121:LYS:CG	2:E:1:ILE:HD13	2.48	0.42
2:E:73:THR:HG23	2:E:75:LYS:NZ	2.33	0.42
2:E:96:ASP:OD1	2:E:96:ASP:C	2.62	0.42
2:B:51:HIS:HA	2:B:65:LEU:O	2.20	0.42
2:B:73:THR:HG23	2:B:75:LYS:NZ	2.35	0.42
1:D:109:PHE:CZ	1:D:111:ARG:HA	2.55	0.42
1:A:48:ARG:HD2	1:A:48:ARG:HA	1.60	0.42
1:A:165:VAL:HG13	1:A:166:GLU:N	2.35	0.42
1:A:230:LEU:HD12	1:A:245:ALA:HB2	2.01	0.42
1:A:253:GLN:NE2	1:A:256:ARG:NE	2.64	0.42
1:A:49:ALA:O	1:A:52:ILE:HG22	2.20	0.41
1:A:182:THR:CG2	4:A:312:HOH:O	2.67	0.41
1:D:103:VAL:HG13	1:D:107:TRP:HA	2.01	0.41
1:A:35:ARG:NE	4:A:327:HOH:O	2.47	0.41
1:D:74:HIS:CE1	1:D:97:ARG:HE	2.38	0.41
1:D:169:ARG:O	1:D:173:GLU:HG3	2.20	0.41
2:E:9:VAL:CG2	2:E:93:VAL:HG22	2.51	0.41
2:E:51:HIS:HA	2:E:65:LEU:O	2.21	0.41
1:A:74:HIS:HD2	1:A:95:VAL:HG11	1.84	0.41
1:A:194:VAL:HG12	1:A:195:SER:N	2.35	0.41
1:D:82:ARG:HG2	1:D:82:ARG:HH11	1.86	0.41
2:E:33:SER:HB3	2:E:62:PHE:CE2	2.56	0.41
1:A:7:TYR:O	1:A:98:MET:HA	2.21	0.40
1:D:45:MET:HE1	3:F:2:LEU:HD11	2.03	0.40
1:A:103:VAL:CG1	1:A:107:TRP:HA	2.48	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:219:ARG:NH1	4:A:319:HOH:O[1_645]	0.97	1.23
1:A:250:PRO:CD	1:D:138:MET:CE[1_456]	2.16	0.04
1:A:138:MET:CE	1:D:250:PRO:CD[1_465]	2.17	0.03

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	274/275 (100%)	268 (98%)	4 (2%)	2 (1%)	18	13
1	D	274/275 (100%)	268 (98%)	5 (2%)	1 (0%)	30	26
2	B	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
2	E	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
3	C	8/10 (80%)	7 (88%)	1 (12%)	0	100	100
3	F	8/10 (80%)	7 (88%)	1 (12%)	0	100	100
All	All	760/770 (99%)	742 (98%)	15 (2%)	3 (0%)	30	26

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	193	ALA
1	D	193	ALA
1	A	194	VAL

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%)	215 (92%)	18 (8%)	12	7
1	D	233/232 (100%)	217 (93%)	16 (7%)	14	9
2	B	95/95 (100%)	86 (90%)	9 (10%)	8	4
2	E	95/95 (100%)	86 (90%)	9 (10%)	8	4

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	10/10 (100%)	9 (90%)	1 (10%)	7	3
3	F	10/10 (100%)	9 (90%)	1 (10%)	7	3
All	All	676/674 (100%)	622 (92%)	54 (8%)	11	7

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	ARG
1	A	48	ARG
1	A	54	GLN
1	A	58	GLU
1	A	103	VAL
1	A	146	LYS
1	A	163	THR
1	A	182	THR
1	A	201	LEU
1	A	212	GLU
1	A	215	LEU
1	A	216	THR
1	A	224	GLN
1	A	225	THR
1	A	228	THR
1	A	249	VAL
1	A	255	GLN
1	A	275	GLU
2	B	1	ILE
2	B	45	ARG
2	B	47	GLU
2	B	48	LYS
2	B	69	GLU
2	B	73	THR
2	B	75	LYS
2	B	85	VAL
2	B	93	VAL
3	C	9	SER
1	D	54	GLN
1	D	89	GLU
1	D	103	VAL
1	D	161	GLU
1	D	163	THR
1	D	165	VAL

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Mol	Chain	Res	Type
1	D	173	GLU
1	D	182	THR
1	D	201	LEU
1	D	212	GLU
1	D	215	LEU
1	D	226	GLN
1	D	228	THR
1	D	249	VAL
1	D	251	SER
1	D	255	GLN
2	E	1	ILE
2	E	45	ARG
2	E	48	LYS
2	E	54	LEU
2	E	73	THR
2	E	74	GLU
2	E	75	LYS
2	E	85	VAL
2	E	93	VAL
3	F	9	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	93	HIS
1	A	114	HIS
1	A	115	GLN
1	A	151	HIS
1	A	192	HIS
1	A	224	GLN
1	A	253	GLN
1	A	263	HIS
1	D	54	GLN
1	D	93	HIS
1	D	114	HIS
1	D	145	HIS
1	D	151	HIS
1	D	155	GLN
1	D	180	GLN
1	D	192	HIS
1	D	218	GLN

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Mol	Chain	Res	Type
1	D	224	GLN
1	D	226	GLN
1	D	253	GLN
1	D	263	HIS
2	E	89	GLN

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	275/275 (100%)	-1.56	0 100 100	6, 14, 33, 51	1 (0%)
1	D	275/275 (100%)	-1.57	0 100 100	6, 15, 33, 51	1 (0%)
2	B	100/100 (100%)	-1.53	0 100 100	5, 18, 38, 53	0
2	E	100/100 (100%)	-1.53	0 100 100	4, 17, 41, 47	0
3	C	10/10 (100%)	-1.65	0 100 100	5, 13, 18, 21	0
3	F	10/10 (100%)	-1.66	0 100 100	5, 14, 20, 21	0
All	All	770/770 (100%)	-1.56	0 100 100	4, 15, 34, 53	2 (0%)

There are no RSRZ outliers to report.

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.