



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:47 AM UTC

PDB ID : 1I1Y / pdb_00001i1y
Title : CRYSTAL STRUCTURE OF HUMAN CLASS I MHC (HLA-A2.1) COM-
PLEXED WITH BETA 2-MICROGLOBULIN AND HIV-RT VARIANT
PEPTIDE I1Y
Authors : Kirksey, T.J.; Collins, E.J.
Deposited on : 1999-04-16
Resolution : 2.20 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

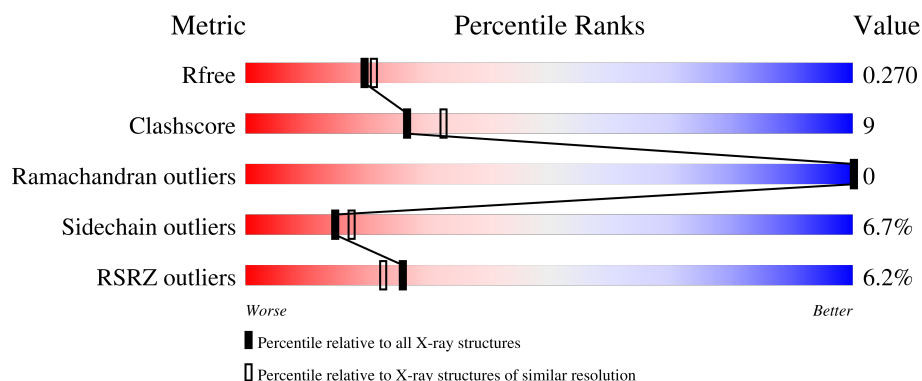
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	8% 76% 19% • •
1	D	275	9% 70% 26% • •
2	B	100	2% 70% 24% 6%
2	E	100	% 67% 26% 6% •
3	C	9	67% 22% 11%

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Mol	Chain	Length	Quality of chain
3	F	9	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: green (67%), yellow (11%), and orange (22%). The percentages are labeled below the bar.

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6515 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 CLASS I HISTOCOMPATIBILITY ANTIGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	0	0
			2247	1403	409	426	9			
1	D	275	Total	C	N	O	S	0	0	0
			2247	1403	409	426	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			

- Molecule 3 is a protein called HIV-RT VARIANT PEPTIDE I1Y (YLKEPVHGV).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			74	49	12	13			
3	F	9	Total	C	N	O	0	0	0
			74	49	12	13			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	64	Total	O	0	0
			64	64		
4	B	23	Total	O	0	0
			23	23		
4	C	5	Total	O	0	0
			5	5		
4	D	67	Total	O	0	0
			67	67		

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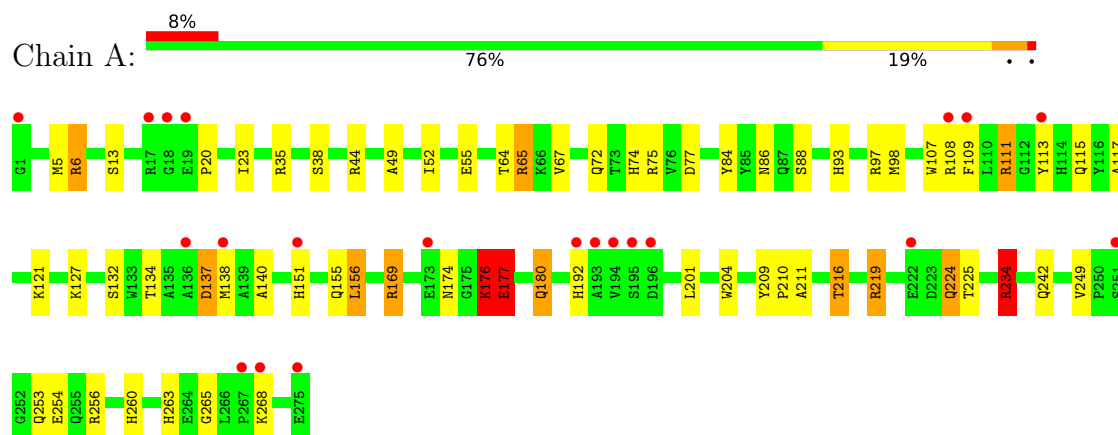
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	37	Total	O	0	0
			37	37		
4	F	3	Total	O	0	0
			3	3		

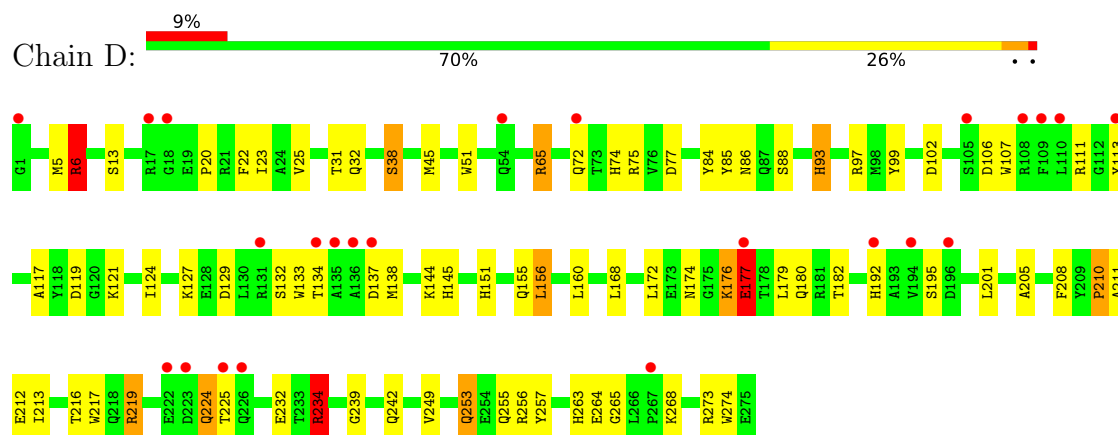
3 Residue-property plots

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.

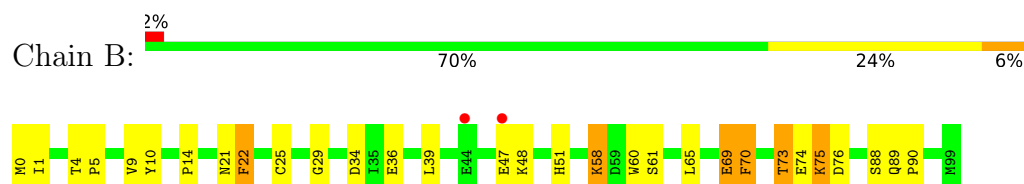
• Molecule 1: CLASS I HISTOCOMPATIBILITY ANTIGEN



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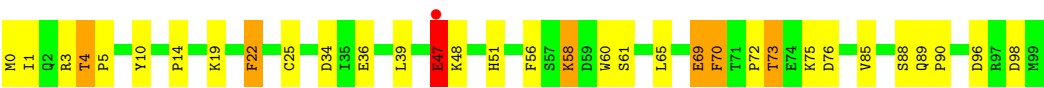


• Molecule 2: BETA 2-MICROGLOBULIN



• Molecule 2: BETA 2-MICROGLOBULIN

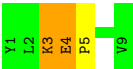




● Molecule 3: HIV-RT VARIANT PEPTIDE I1Y (YLKEPVHGV)



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	50.61Å 63.59Å 75.33Å 81.93° 75.94° 78.00°	Depositor
Resolution (Å)	15.00 – 2.20 15.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.0 (15.00-2.20) 97.6 (15.00-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.82 (at 2.20Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.233 , 0.289 0.231 , 0.270	Depositor DCC
R_{free} test set	3615 reflections (8.20%)	wwPDB-VP
Wilson B-factor (Å ²)	19.0	Xtriage
Anisotropy	0.398	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 50.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6515	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.17% 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	0.71	0/2312	1.81	38/3137 (1.2%)
1	D	0.76	2/2312 (0.1%)	1.81	45/3137 (1.4%)
2	B	0.67	0/860	1.57	11/1162 (0.9%)
2	E	0.68	0/860	1.55	15/1162 (1.3%)
3	C	0.76	0/76	1.21	0/101
3	F	0.83	0/76	1.29	1/101 (1.0%)
All	All	0.72	2/6496 (0.0%)	1.74	110/8800 (1.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	217	TRP	NE1-CE2	-5.60	1.31	1.37
1	D	133	TRP	NE1-CE2	-5.22	1.31	1.37

All (110) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	234	ARG	CD-NE-CZ	38.39	178.15	124.40
1	D	234	ARG	CD-NE-CZ	33.98	171.97	124.40
1	A	224	GLN	OE1-CD-NE2	-18.27	104.33	122.60
1	D	224	GLN	OE1-CD-NE2	-15.34	107.26	122.60
1	D	65	ARG	NE-CZ-NH2	12.73	130.66	119.20
1	A	65	ARG	NE-CZ-NH2	11.10	129.19	119.20
1	D	65	ARG	NE-CZ-NH1	-10.38	111.12	121.50
1	A	219	ARG	CD-NE-CZ	9.56	137.79	124.40
1	A	65	ARG	NE-CZ-NH1	-9.42	112.08	121.50
1	D	219	ARG	CD-NE-CZ	9.38	137.54	124.40
1	A	224	GLN	CG-CD-OE1	8.97	138.74	120.80
1	D	6	ARG	CD-NE-CZ	8.62	136.47	124.40
1	A	6	ARG	NE-CZ-NH1	8.36	129.86	121.50
1	A	234	ARG	NE-CZ-NH2	8.14	126.52	119.20
1	A	209	TYR	CA-C-O	-8.01	111.06	119.22
1	D	224	GLN	CG-CD-OE1	7.71	136.23	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	74	GLU	CA-C-N	7.69	131.09	120.63
2	B	74	GLU	C-N-CA	7.69	131.09	120.63
2	B	34	ASP	CA-CB-CG	7.67	120.28	112.60
1	D	151	HIS	CA-CB-CG	-7.38	106.42	113.80
1	A	6	ARG	CD-NE-CZ	7.38	134.74	124.40
1	D	213	ILE	CA-C-O	-7.22	115.82	120.66
1	A	108	ARG	NE-CZ-NH1	7.17	128.67	121.50
2	B	70	PHE	CA-CB-CG	7.16	120.96	113.80
1	A	151	HIS	CA-CB-CG	-7.09	106.71	113.80
1	A	137	ASP	CA-CB-CG	-7.06	105.54	112.60
2	E	14	PRO	CA-C-O	-6.99	113.70	121.32
1	A	5	MET	CA-C-O	-6.89	112.93	120.30
1	D	224	GLN	CA-C-O	6.79	128.16	119.95
1	D	6	ARG	NE-CZ-NH1	6.60	128.10	121.50
1	D	192	HIS	CA-CB-CG	6.51	120.31	113.80
1	D	86	ASN	OD1-CG-ND2	-6.45	116.15	122.60
1	A	174	ASN	OD1-CG-ND2	-6.44	116.16	122.60
1	D	211	ALA	N-CA-C	6.38	118.23	111.28
1	D	5	MET	CA-C-O	-6.37	113.48	120.36
1	D	253	GLN	OE1-CD-NE2	-6.34	116.25	122.60
1	A	260	HIS	CA-CB-CG	-6.33	107.47	113.80
1	D	25	VAL	O-C-N	6.26	130.00	123.05
1	D	23	ILE	CA-C-O	-6.19	113.89	120.39
1	A	109	PHE	CA-C-O	-6.19	115.07	120.88
2	B	21	ASN	CA-CB-CG	6.17	118.78	112.60
1	D	219	ARG	NE-CZ-NH1	6.17	127.67	121.50
1	D	210	PRO	CA-C-N	6.17	128.54	120.28
1	D	210	PRO	C-N-CA	6.17	128.54	120.28
1	D	232	GLU	O-C-N	-6.11	115.44	122.89
1	D	234	ARG	CG-CD-NE	6.11	125.43	112.00
2	E	98	ASP	CA-CB-CG	6.06	118.66	112.60
1	D	177	GLU	CB-CG-CD	6.06	122.90	112.60
2	E	3	ARG	NE-CZ-NH2	6.01	124.61	119.20
1	A	180	GLN	OE1-CD-NE2	-5.99	116.61	122.60
2	B	75	LYS	N-CA-C	5.95	117.63	111.03
1	D	211	ALA	CA-C-O	-5.90	114.30	120.55
1	A	23	ILE	CB-CA-C	-5.88	101.93	110.63
1	A	111	ARG	CA-C-N	5.87	126.13	121.61
1	A	111	ARG	C-N-CA	5.87	126.13	121.61
1	D	93	HIS	CA-CB-CG	-5.87	107.93	113.80
1	A	176	LYS	CA-C-N	5.86	128.14	120.28
1	A	176	LYS	C-N-CA	5.86	128.14	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	25	VAL	N-CA-CB	5.84	121.03	111.45
1	D	182	THR	N-CA-CB	5.83	120.88	110.80
1	D	32	GLN	CA-C-O	-5.83	114.64	121.16
1	D	84	TYR	CA-CB-CG	5.78	124.30	113.90
2	E	69	GLU	CB-CG-CD	5.77	122.41	112.60
2	E	22	PHE	CA-CB-CG	5.74	119.54	113.80
1	D	274	TRP	CA-C-O	-5.72	114.13	120.70
2	E	96	ASP	O-C-N	-5.72	116.52	123.27
1	A	84	TYR	CA-CB-CG	5.70	124.16	113.90
2	E	4	THR	CB-CA-C	5.68	116.89	108.87
2	B	5	PRO	CB-CA-C	-5.67	104.50	111.64
1	A	65	ARG	CD-NE-CZ	-5.65	116.49	124.40
2	E	70	PHE	CA-CB-CG	5.65	119.45	113.80
1	D	23	ILE	CB-CA-C	-5.64	102.28	110.63
1	A	192	HIS	CA-CB-CG	5.62	119.42	113.80
1	D	212	GLU	CA-C-O	-5.59	115.04	121.19
1	D	239	GLY	N-CA-C	-5.59	107.65	114.92
1	D	160	LEU	CA-C-O	-5.59	113.55	119.97
1	A	77	ASP	CA-CB-CG	5.57	118.17	112.60
1	A	86	ASN	OD1-CG-ND2	-5.56	117.04	122.60
2	E	4	THR	CA-C-O	5.56	124.86	120.19
1	D	205	ALA	N-CA-CB	5.54	119.47	110.77
1	A	55	GLU	CA-C-O	-5.50	115.57	121.56
1	D	31	THR	O-C-N	-5.46	116.99	123.22
2	E	61	SER	CB-CA-C	5.43	118.93	109.65
2	E	3	ARG	NH1-CZ-NH2	-5.42	112.25	119.30
2	B	14	PRO	CA-C-O	-5.41	115.42	121.32
1	D	77	ASP	CA-C-O	-5.41	114.79	120.63
1	D	208	PHE	CA-CB-CG	5.40	119.20	113.80
1	D	99	TYR	CA-CB-CG	5.39	123.60	113.90
1	D	86	ASN	CA-CB-CG	5.38	117.98	112.60
1	A	177	GLU	CB-CG-CD	5.33	121.66	112.60
1	A	115	GLN	OE1-CD-NE2	-5.30	117.30	122.60
2	E	5	PRO	CB-CA-C	-5.27	104.20	111.21
2	E	34	ASP	CA-CB-CG	5.27	117.87	112.60
1	A	86	ASN	CA-CB-CG	5.26	117.86	112.60
1	A	211	ALA	N-CA-C	5.23	117.72	111.71
2	E	56	PHE	CA-CB-CG	5.22	119.03	113.80
1	A	169	ARG	NE-CZ-NH1	-5.22	116.28	121.50
2	B	9	VAL	CB-CA-C	-5.21	103.35	110.96
1	D	257	TYR	N-CA-C	5.18	118.26	109.76
1	A	204	TRP	N-CA-C	5.14	117.87	109.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	4	GLU	O-C-N	-5.13	117.84	121.84
2	B	22	PHE	CA-CB-CG	5.12	118.92	113.80
1	A	35	ARG	O-C-N	5.11	128.89	123.42
1	D	45	MET	CA-C-O	-5.10	115.45	121.16
2	B	69	GLU	CB-CG-CD	5.08	121.24	112.60
1	A	174	ASN	CA-CB-CG	5.04	117.64	112.60
1	D	129	ASP	CA-CB-CG	5.03	117.63	112.60
2	E	47	GLU	CB-CG-CD	5.03	121.15	112.60
1	A	77	ASP	CA-C-O	-5.03	115.22	120.55
1	D	174	ASN	CA-CB-CG	5.01	117.61	112.60

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	2247	0	2096	43	0
1	D	2247	0	2096	47	0
2	B	837	0	803	19	0
2	E	837	0	803	22	0
3	C	74	0	76	8	0
3	F	74	0	76	7	0
4	A	64	0	0	2	0
4	B	23	0	0	0	0
4	C	5	0	0	0	0
4	D	67	0	0	7	0
4	E	37	0	0	3	0
4	F	3	0	0	0	0
All	All	6515	0	5950	114	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 9.

All (114) 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:93:HIS:CD2	2:B:0:MET:HE1	1.64	1.31
1:A:93:HIS:NE2	2:B:0:MET:HE1	1.49	1.25
1:D:93:HIS:CD2	2:E:0:MET:HE1	1.75	1.20
1:D:93:HIS:NE2	2:E:0:MET:HE1	1.55	1.18
1:D:93:HIS:CD2	2:E:0:MET:CE	2.34	1.10
1:A:93:HIS:CD2	2:B:0:MET:CE	2.37	1.06
1:D:93:HIS:NE2	2:E:0:MET:CE	2.24	1.00
2:E:47:GLU:HG2	4:E:102:HOH:O	1.60	0.99
1:A:93:HIS:NE2	2:B:0:MET:CE	2.30	0.90
1:D:234:ARG:HD2	1:D:242:GLN:HB2	1.54	0.88
1:A:234:ARG:HD2	1:A:242:GLN:HB2	1.57	0.86
1:D:72:GLN:HE22	1:D:75:ARG:HH11	1.27	0.82
1:D:93:HIS:CD2	2:E:0:MET:HE2	2.16	0.81
1:A:155:GLN:HE21	3:C:3:LYS:NZ	1.83	0.76
1:A:155:GLN:HE21	3:C:3:LYS:HZ3	1.30	0.76
1:A:72:GLN:HE22	1:A:75:ARG:HH11	1.32	0.75
2:E:89:GLN:HB2	2:E:90:PRO:HD2	1.70	0.74
2:E:73:THR:HG22	2:E:76:ASP:H	1.56	0.71
3:C:3:LYS:HG3	3:C:4:GLU:N	2.07	0.69
2:B:89:GLN:HB2	2:B:90:PRO:HD2	1.75	0.69
2:B:58:LYS:HD2	2:B:58:LYS:H	1.58	0.68
1:D:155:GLN:NE2	3:F:5:PRO:HD2	2.09	0.68
1:D:263:HIS:HD2	1:D:265:GLY:H	1.41	0.67
2:E:58:LYS:H	2:E:58:LYS:HD2	1.59	0.67
1:D:201:LEU:HD22	1:D:249:VAL:HG21	1.75	0.67
2:B:73:THR:HG22	2:B:76:ASP:H	1.59	0.67
1:D:65:ARG:NH1	4:D:303:HOH:O	2.27	0.65
1:D:263:HIS:CD2	1:D:265:GLY:H	2.15	0.65
1:A:65:ARG:NH1	4:A:329:HOH:O	2.29	0.65
1:A:93:HIS:HD2	2:B:0:MET:CE	2.03	0.65
2:E:75:LYS:HA	4:E:127:HOH:O	1.97	0.64
1:A:263:HIS:CD2	1:A:265:GLY:H	2.17	0.62
1:A:93:HIS:CD2	2:B:0:MET:HE2	2.34	0.62
1:D:155:GLN:HE21	3:F:3:LYS:NZ	1.99	0.60
1:D:219:ARG:HB2	1:D:224:GLN:HE21	1.65	0.60
1:D:176:LYS:HG2	1:D:180:GLN:NE2	2.16	0.60
3:F:3:LYS:HG3	3:F:4:GLU:N	2.14	0.60
1:D:93:HIS:HD2	2:E:0:MET:HE2	1.64	0.59
1:A:155:GLN:NE2	3:C:5:PRO:HD2	2.18	0.58
1:D:177:GLU:H	1:D:177:GLU:CD	2.12	0.58
1:A:176:LYS:HG2	1:A:180:GLN:NE2	2.19	0.58
1:A:263:HIS:HD2	1:A:265:GLY:H	1.50	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:ARG:HB3	1:A:224:GLN:HE21	1.70	0.57
1:D:155:GLN:HE21	3:F:3:LYS:HZ3	1.53	0.57
1:A:201:LEU:HD22	1:A:249:VAL:HG21	1.88	0.56
2:E:85:VAL:HG13	4:E:100:HOH:O	2.06	0.55
1:A:253:GLN:HE21	1:A:256:ARG:HD3	1.72	0.55
1:D:127:LYS:HE3	1:D:134:THR:OG1	2.06	0.55
1:A:216:THR:HG21	4:A:296:HOH:O	2.07	0.54
1:D:74:HIS:CE1	1:D:97:ARG:HE	2.26	0.54
1:A:74:HIS:CE1	1:A:97:ARG:HE	2.26	0.54
1:D:93:HIS:HD2	2:E:0:MET:CE	2.11	0.54
1:A:219:ARG:CB	1:A:224:GLN:HE21	2.21	0.53
1:D:234:ARG:HD3	2:E:10:TYR:CE2	2.42	0.53
1:D:72:GLN:NE2	1:D:75:ARG:HH11	2.02	0.53
1:A:121:LYS:HG3	2:B:1:ILE:HD12	1.89	0.52
2:E:25:CYS:HB2	2:E:39:LEU:HD21	1.91	0.52
1:A:177:GLU:H	1:A:177:GLU:CD	2.17	0.51
1:A:253:GLN:NE2	1:A:256:ARG:HH11	2.09	0.50
2:E:89:GLN:HB2	2:E:90:PRO:CD	2.40	0.50
2:B:25:CYS:HB2	2:B:39:LEU:HD21	1.93	0.50
1:D:264:GLU:HB2	4:D:305:HOH:O	2.11	0.50
1:A:107:TRP:O	1:A:169:ARG:NH2	2.44	0.50
1:A:121:LYS:HG3	2:B:1:ILE:CD1	2.42	0.50
1:D:6:ARG:NH2	1:D:102:ASP:OD1	2.45	0.50
3:C:3:LYS:CG	3:C:4:GLU:N	2.74	0.49
1:D:156:LEU:HD12	3:F:3:LYS:HE3	1.94	0.49
1:D:255:GLN:HG2	4:D:321:HOH:O	2.12	0.49
1:A:72:GLN:NE2	1:A:75:ARG:HH11	2.05	0.48
1:A:155:GLN:NE2	3:C:3:LYS:NZ	2.58	0.48
1:A:127:LYS:HE3	1:A:134:THR:OG1	2.14	0.47
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.49	0.47
1:D:111:ARG:HG3	1:D:113:TYR:CZ	2.49	0.47
1:D:124:ILE:HG13	1:D:134:THR:O	2.15	0.47
2:E:51:HIS:HA	2:E:65:LEU:O	2.14	0.47
1:A:155:GLN:NE2	3:C:3:LYS:HZ3	2.05	0.47
3:F:3:LYS:CG	3:F:4:GLU:N	2.77	0.47
1:A:93:HIS:HD2	2:B:0:MET:HE2	1.75	0.47
1:D:6:ARG:HD2	4:D:293:HOH:O	2.14	0.47
1:D:144:LYS:HE2	4:D:297:HOH:O	2.14	0.46
2:B:29:GLY:HA2	2:B:61:SER:OG	2.16	0.46
1:D:22:PHE:H	1:D:38:SER:HB2	1.81	0.46
1:A:49:ALA:O	1:A:52:ILE:HG22	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:PRO:O	1:A:263:HIS:HE1	1.99	0.45
1:D:117:ALA:HB2	2:E:60:TRP:CE2	2.51	0.45
1:A:234:ARG:HD3	2:B:10:TYR:CE2	2.53	0.44
2:E:19:LYS:O	2:E:72:PRO:HD2	2.18	0.44
1:D:253:GLN:NE2	1:D:256:ARG:HH11	2.15	0.44
1:A:65:ARG:HH11	1:A:65:ARG:HD3	1.34	0.44
2:B:51:HIS:HA	2:B:65:LEU:O	2.17	0.44
1:A:111:ARG:HG3	1:A:113:TYR:CZ	2.52	0.44
1:A:137:ASP:H	1:A:140:ALA:HB3	1.83	0.43
1:D:156:LEU:CD1	3:F:3:LYS:HE3	2.47	0.43
2:B:58:LYS:H	2:B:58:LYS:CD	2.29	0.43
2:E:22:PHE:CE2	2:E:69:GLU:HG3	2.53	0.43
1:D:253:GLN:HE21	1:D:256:ARG:HD3	1.83	0.43
1:D:65:ARG:HH11	1:D:65:ARG:HD3	1.36	0.43
1:A:44:ARG:HA	1:A:64:THR:HG23	2.00	0.43
1:D:85:TYR:OH	1:D:137:ASP:OD2	2.17	0.42
1:A:156:LEU:HD12	3:C:3:LYS:HE3	2.00	0.42
2:B:22:PHE:CE2	2:B:69:GLU:HG3	2.55	0.42
1:D:13:SER:HA	1:D:20:PRO:HB3	2.02	0.42
1:D:210:PRO:O	1:D:263:HIS:HE1	2.02	0.42
1:D:273:ARG:NH1	4:D:333:HOH:O	2.52	0.42
1:D:51:TRP:CZ2	1:D:179:LEU:HD11	2.55	0.41
1:D:168:LEU:O	1:D:172:LEU:HG	2.20	0.41
1:A:98:MET:SD	1:A:98:MET:C	3.03	0.41
1:A:13:SER:HA	1:A:20:PRO:HB3	2.02	0.41
1:D:93:HIS:HD2	1:D:119:ASP:OD2	2.04	0.41
1:A:253:GLN:O	1:A:254:GLU:C	2.62	0.41
1:D:121:LYS:HG3	2:E:1:ILE:HD12	2.03	0.40
1:D:145:HIS:HD2	4:D:280:HOH:O	2.04	0.40
2:E:73:THR:HG23	2:E:75:LYS:H	1.87	0.40
1:D:106:ASP:O	1:D:107:TRP:HB2	2.22	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	273/275 (99%)	265 (97%)	8 (3%)	0	100	100
1	D	273/275 (99%)	264 (97%)	9 (3%)	0	100	100
2	B	98/100 (98%)	95 (97%)	3 (3%)	0	100	100
2	E	98/100 (98%)	96 (98%)	2 (2%)	0	100	100
3	C	7/9 (78%)	7 (100%)	0	0	100	100
3	F	7/9 (78%)	7 (100%)	0	0	100	100
All	All	756/768 (98%)	734 (97%)	22 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	231/231 (100%)	218 (94%)	13 (6%)	19	24
1	D	231/231 (100%)	218 (94%)	13 (6%)	19	24
2	B	95/95 (100%)	86 (90%)	9 (10%)	8	8
2	E	95/95 (100%)	87 (92%)	8 (8%)	10	11
3	C	8/8 (100%)	7 (88%)	1 (12%)	4	4
3	F	8/8 (100%)	7 (88%)	1 (12%)	4	4
All	All	668/668 (100%)	623 (93%)	45 (7%)	15	17

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
1	A	38	SER
1	A	67	VAL
1	A	88	SER

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Mol	Chain	Res	Type
1	A	132	SER
1	A	138	MET
1	A	156	LEU
1	A	176	LYS
1	A	177	GLU
1	A	216	THR
1	A	225	THR
1	A	234	ARG
1	A	268	LYS
2	B	4	THR
2	B	36	GLU
2	B	47	GLU
2	B	48	LYS
2	B	58	LYS
2	B	70	PHE
2	B	73	THR
2	B	75	LYS
2	B	88	SER
3	C	3	LYS
1	D	6	ARG
1	D	38	SER
1	D	88	SER
1	D	132	SER
1	D	138	MET
1	D	156	LEU
1	D	176	LYS
1	D	177	GLU
1	D	195	SER
1	D	216	THR
1	D	225	THR
1	D	234	ARG
1	D	268	LYS
2	E	4	THR
2	E	36	GLU
2	E	47	GLU
2	E	48	LYS
2	E	58	LYS
2	E	70	PHE
2	E	73	THR
2	E	88	SER
3	F	3	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	HIS
1	A	72	GLN
1	A	74	HIS
1	A	93	HIS
1	A	155	GLN
1	A	174	ASN
1	A	180	GLN
1	A	192	HIS
1	A	224	GLN
1	A	253	GLN
1	A	263	HIS
2	B	2	GLN
2	B	83	ASN
2	B	89	GLN
3	C	7	HIS
1	D	70	HIS
1	D	72	GLN
1	D	74	HIS
1	D	93	HIS
1	D	145	HIS
1	D	155	GLN
1	D	224	GLN
1	D	253	GLN
1	D	263	HIS
2	E	2	GLN
2	E	83	ASN
2	E	89	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	275/275 (100%)	0.39	21 (7%) 20 17	9, 21, 40, 57	0
1	D	275/275 (100%)	0.37	24 (8%) 16 13	9, 21, 40, 57	0
2	B	100/100 (100%)	-0.04	2 (2%) 65 62	9, 17, 34, 38	0
2	E	100/100 (100%)	0.00	1 (1%) 79 77	9, 17, 34, 38	0
3	C	9/9 (100%)	-0.08	0 100 100	14, 19, 24, 29	0
3	F	9/9 (100%)	-0.15	0 100 100	14, 20, 24, 29	0
All	All	768/768 (100%)	0.26	48 (6%) 26 23	9, 20, 38, 57	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	105	SER	4.7
1	D	192	HIS	4.1
1	D	196	ASP	4.0
2	E	47	GLU	3.8
1	A	196	ASP	3.8
1	D	17	ARG	3.8
1	A	108	ARG	3.6
1	A	1	GLY	3.5
1	D	225	THR	3.4
1	A	113	TYR	3.4
1	D	223	ASP	3.4
2	B	44	GLU	3.3
1	A	193	ALA	3.2
1	A	136	ALA	2.8
1	A	192	HIS	2.8
1	A	267	PRO	2.7
1	D	136	ALA	2.7
1	D	1	GLY	2.7
1	D	54	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	226	GLN	2.6
1	A	195	SER	2.6
1	A	109	PHE	2.5
1	D	109	PHE	2.5
1	D	222	GLU	2.5
1	D	108	ARG	2.5
1	A	173	GLU	2.4
1	A	18	GLY	2.4
2	B	47	GLU	2.4
1	A	17	ARG	2.4
1	D	72	GLN	2.4
1	D	113	TYR	2.3
1	D	137	ASP	2.3
1	D	194	VAL	2.3
1	A	268	LYS	2.2
1	A	151	HIS	2.2
1	D	134	THR	2.2
1	D	131	ARG	2.2
1	A	19	GLU	2.2
1	A	275	GLU	2.2
1	D	177	GLU	2.2
1	A	251	SER	2.2
1	D	135	ALA	2.2
1	A	222	GLU	2.1
1	A	138	MET	2.1
1	D	267	PRO	2.1
1	D	18	GLY	2.1
1	A	194	VAL	2.0
1	D	110	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.