



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:42 PM UTC

PDB ID : 1HHK / pdb_00001hhk
Title : THE ANTIGENIC IDENTITY OF PEPTIDE(SLASH)MHC COMPLEXES:
A COMPARISON OF THE CONFORMATION OF FIVE PEPTIDES PRE-
SENTED BY HLA-A2
Authors : Madden, D.R.; Garboczi, D.N.; Wiley, D.C.
Deposited on : 1993-06-30
Resolution : 2.50 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

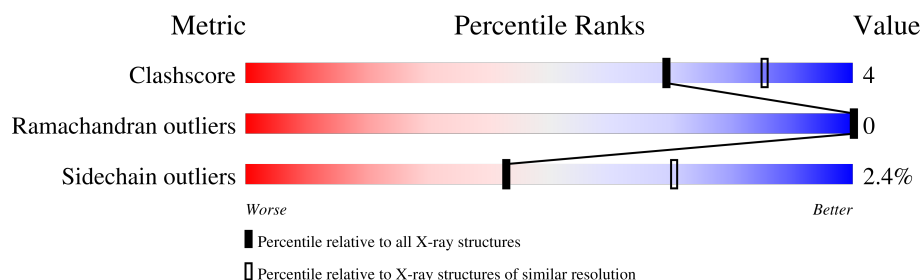
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	275	76% 21% ..
1	D	275	80% 17% ..
2	B	100	73% 25% .
2	E	100	78% 20% .
3	C	9	56% 44%
3	F	9	56% 44%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6322 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 (HLA-A*0201) (ALPHA CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	14	0	0
			2247	1403	409	426	9			
1	D	275	Total	C	N	O	S	14	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	10	0	0
			837	533	141	159	4			
2	E	100	Total	C	N	O	S	10	0	0
			837	533	141	159	4			

- Molecule 3 is a protein called NONAMERIC PEPTIDE FROM HTLV-1 TAX PROTEIN (RESIDUES 11-19).

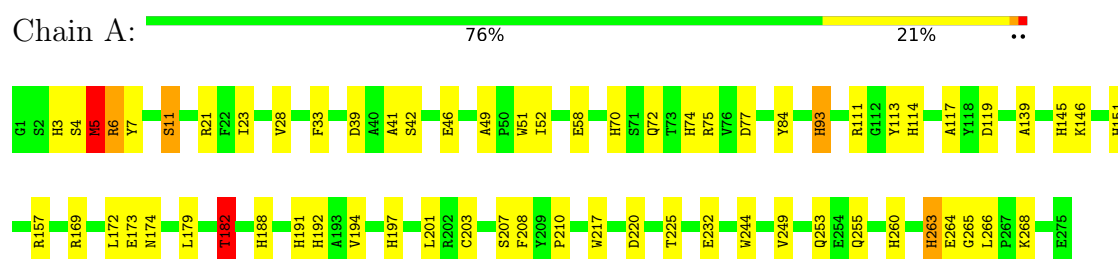
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			77	56	9	12			
3	F	9	Total	C	N	O	0	0	0
			77	56	9	12			

3 Residue-property plots

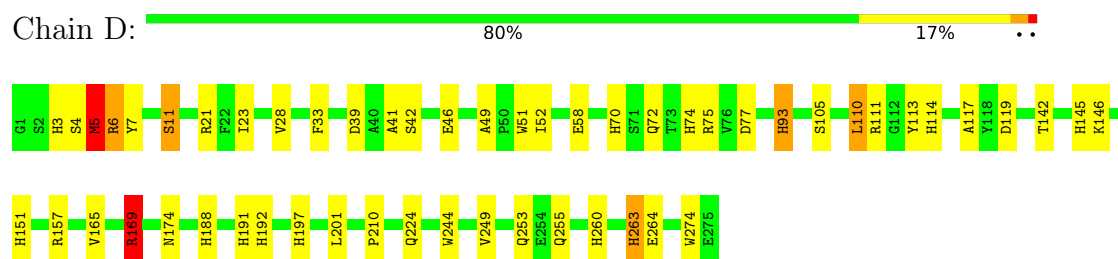
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

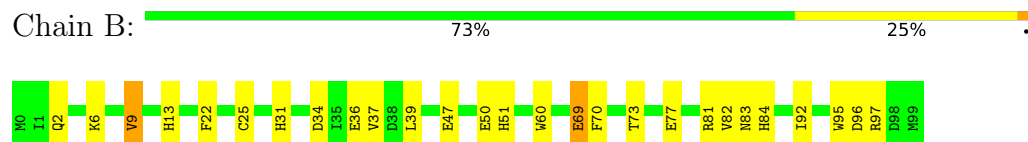
• Molecule 1: CLASS I HISTOCOMPATIBILITY ANTIGEN (HLA-A*0201) (ALPHA CHAIN)



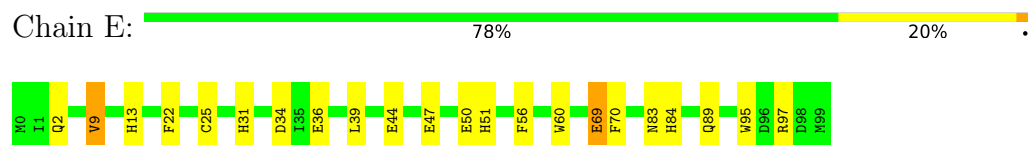
• Molecule 1: CLASS I HISTOCOMPATIBILITY ANTIGEN (HLA-A*0201) (ALPHA CHAIN)



• Molecule 2: BETA 2-MICROGLOBULIN



• Molecule 2: BETA 2-MICROGLOBULIN



• Molecule 3: NONAMERIC PEPTIDE FROM HTLV-1 TAX PROTEIN (RESIDUES 11-19)





- Molecule 3: NONAMERIC PEPTIDE FROM HTLV-1 TAX PROTEIN (RESIDUES 11-19)

Chain F:  56% 44%

A horizontal bar chart for Chain F. The bar is divided into two segments: a green segment on the left representing 56% and a yellow segment on the right representing 44%.



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	50.56Å 63.79Å 75.08Å 81.58° 75.66° 77.38°	Depositor
Resolution (Å)	(Not available) – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.262 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6322	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

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	1.05	15/2312 (0.6%)	1.70	45/3137 (1.4%)
1	D	1.06	16/2312 (0.7%)	1.69	36/3137 (1.1%)
2	B	1.03	6/860 (0.7%)	1.59	13/1162 (1.1%)
2	E	1.05	5/860 (0.6%)	1.62	15/1162 (1.3%)
3	C	0.66	0/80	1.44	1/108 (0.9%)
3	F	0.73	0/80	1.47	1/108 (0.9%)
All	All	1.04	42/6504 (0.6%)	1.67	111/8814 (1.3%)

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	93	HIS	CD2-NE2	-7.45	1.29	1.37
1	D	260	HIS	CD2-NE2	-7.38	1.29	1.37
1	A	260	HIS	CD2-NE2	-7.33	1.29	1.37
1	D	188	HIS	CD2-NE2	-7.09	1.30	1.37
2	E	84	HIS	CD2-NE2	-7.04	1.30	1.37
1	A	191	HIS	CD2-NE2	-6.96	1.30	1.37
2	E	31	HIS	CD2-NE2	-6.94	1.30	1.37
2	B	31	HIS	CD2-NE2	-6.93	1.30	1.37
2	B	84	HIS	CD2-NE2	-6.93	1.30	1.37
1	A	151	HIS	CD2-NE2	-6.89	1.30	1.37
1	D	93	HIS	CD2-NE2	-6.85	1.30	1.37
2	E	13	HIS	CD2-NE2	-6.78	1.30	1.37
1	D	191	HIS	CD2-NE2	-6.63	1.30	1.37
1	A	263	HIS	CD2-NE2	-6.58	1.30	1.37
1	D	263	HIS	CD2-NE2	-6.52	1.30	1.37
1	A	114	HIS	CD2-NE2	-6.50	1.30	1.37
1	D	3	HIS	CG-ND1	-6.38	1.31	1.38
1	D	145	HIS	CD2-NE2	-6.38	1.30	1.37
1	A	74	HIS	CD2-NE2	-6.35	1.30	1.37
2	B	13	HIS	CD2-NE2	-6.33	1.30	1.37
1	A	188	HIS	CD2-NE2	-6.27	1.30	1.37
1	D	197	HIS	CD2-NE2	-6.24	1.30	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	151	HIS	CD2-NE2	-6.22	1.31	1.37
1	A	197	HIS	CD2-NE2	-6.14	1.31	1.37
1	D	114	HIS	CD2-NE2	-6.00	1.31	1.37
1	D	3	HIS	CD2-NE2	-5.98	1.31	1.37
1	A	3	HIS	CG-ND1	-5.95	1.31	1.38
1	A	70	HIS	CD2-NE2	-5.93	1.31	1.37
1	D	74	HIS	CD2-NE2	-5.92	1.31	1.37
2	B	51	HIS	CD2-NE2	-5.83	1.31	1.37
1	D	70	HIS	CD2-NE2	-5.74	1.31	1.37
1	A	145	HIS	CD2-NE2	-5.64	1.31	1.37
1	D	192	HIS	CD2-NE2	-5.59	1.31	1.37
1	A	151	HIS	CG-ND1	-5.51	1.32	1.38
2	E	51	HIS	CD2-NE2	-5.42	1.31	1.37
1	A	192	HIS	CD2-NE2	-5.39	1.31	1.37
2	B	9	VAL	CA-CB	5.34	1.60	1.54
1	A	3	HIS	CD2-NE2	-5.29	1.32	1.37
1	D	188	HIS	CG-ND1	-5.27	1.32	1.38
1	D	191	HIS	CG-ND1	-5.18	1.32	1.38
2	B	51	HIS	CG-ND1	-5.11	1.32	1.38
2	E	9	VAL	CA-CB	5.08	1.60	1.54

All (111) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	5	MET	CG-SD-CE	-11.57	75.44	100.90
1	D	5	MET	CG-SD-CE	-10.61	77.57	100.90
1	D	174	ASN	OD1-CG-ND2	-8.57	114.03	122.60
1	A	174	ASN	OD1-CG-ND2	-8.33	114.27	122.60
2	E	89	GLN	CA-CB-CG	-8.22	97.66	114.10
2	B	47	GLU	N-CA-C	7.75	119.80	111.36
1	A	194	VAL	N-CA-C	-7.68	104.31	111.45
1	D	260	HIS	CB-CG-CD2	-7.22	121.81	131.20
2	B	31	HIS	CB-CG-CD2	-7.22	121.82	131.20
1	D	263	HIS	CB-CG-CD2	-6.98	122.12	131.20
1	D	169	ARG	CA-CB-CG	6.98	128.06	114.10
1	A	33	PHE	CA-CB-CG	6.95	120.75	113.80
1	D	93	HIS	CB-CG-CD2	-6.92	122.20	131.20
2	E	31	HIS	CB-CG-CD2	-6.92	122.21	131.20
1	A	260	HIS	CB-CG-CD2	-6.89	122.25	131.20
1	D	110	LEU	N-CA-C	-6.89	104.75	113.43
1	A	93	HIS	CB-CG-CD2	-6.72	122.47	131.20
2	E	47	GLU	N-CA-C	6.67	118.35	111.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	84	HIS	CB-CG-CD2	-6.52	122.72	131.20
3	F	7	VAL	N-CA-C	-6.48	98.70	108.17
2	B	96	ASP	N-CA-C	-6.40	99.29	109.59
2	B	77	GLU	CA-CB-CG	6.38	126.86	114.10
2	E	2	GLN	OE1-CD-NE2	-6.37	116.23	122.60
3	C	7	VAL	N-CA-C	-6.35	98.89	108.17
2	B	84	HIS	CB-CG-CD2	-6.33	122.97	131.20
1	A	263	HIS	CB-CG-CD2	-6.31	123.00	131.20
1	A	191	HIS	CB-CG-CD2	-6.25	123.07	131.20
1	D	33	PHE	CA-CB-CG	6.24	120.04	113.80
1	A	253	GLN	OE1-CD-NE2	-6.17	116.43	122.60
1	D	157	ARG	CG-CD-NE	-6.17	98.42	112.00
2	B	69	GLU	CB-CG-CD	6.12	123.01	112.60
2	B	2	GLN	OE1-CD-NE2	-6.12	116.48	122.60
1	D	46	GLU	CA-C-N	6.12	125.88	119.76
1	D	46	GLU	C-N-CA	6.12	125.88	119.76
1	A	191	HIS	CA-CB-CG	6.09	119.89	113.80
1	D	145	HIS	CB-CG-CD2	-6.01	123.39	131.20
1	A	72	GLN	CA-C-O	-5.99	114.07	120.42
1	A	28	VAL	N-CA-C	-5.98	99.02	107.75
1	D	72	GLN	CA-C-O	-5.97	114.09	120.42
2	E	97	ARG	CA-CB-CG	5.96	126.02	114.10
2	E	69	GLU	CB-CG-CD	5.95	122.72	112.60
1	D	197	HIS	CB-CG-CD2	-5.95	123.47	131.20
1	D	253	GLN	OE1-CD-NE2	-5.95	116.65	122.60
1	D	28	VAL	N-CA-C	-5.93	99.09	107.75
1	D	191	HIS	CB-CG-CD2	-5.90	123.53	131.20
2	E	44	GLU	CB-CG-CD	5.86	122.56	112.60
2	B	13	HIS	CB-CG-CD2	-5.84	123.60	131.20
1	A	93	HIS	CB-CG-ND1	5.82	131.43	122.70
1	A	6	ARG	CA-CB-CG	-5.82	102.47	114.10
1	A	197	HIS	CB-CG-CD2	-5.80	123.67	131.20
2	E	56	PHE	CA-CB-CG	5.79	119.59	113.80
1	A	264	GLU	CA-CB-CG	5.78	125.65	114.10
1	A	151	HIS	CB-CG-CD2	-5.77	123.70	131.20
1	A	3	HIS	CA-CB-CG	-5.71	108.09	113.80
1	D	77	ASP	CA-CB-CG	5.70	118.30	112.60
1	A	188	HIS	CB-CG-CD2	-5.66	123.84	131.20
1	A	255	GLN	OE1-CD-NE2	-5.66	116.94	122.60
1	A	41	ALA	N-CA-C	5.65	117.44	111.28
1	A	46	GLU	CA-C-N	5.65	125.41	119.76
1	A	46	GLU	C-N-CA	5.65	125.41	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	6	ARG	CA-CB-CG	-5.65	102.80	114.10
1	D	188	HIS	CB-CG-CD2	-5.64	123.87	131.20
1	A	157	ARG	CG-CD-NE	-5.63	99.61	112.00
1	D	93	HIS	CB-CG-ND1	5.62	131.13	122.70
1	D	264	GLU	CA-CB-CG	5.61	125.32	114.10
2	E	13	HIS	CB-CG-CD2	-5.60	123.92	131.20
1	A	77	ASP	CA-CB-CG	5.55	118.15	112.60
2	B	70	PHE	CA-CB-CG	5.49	119.29	113.80
2	B	13	HIS	CB-CG-ND1	5.47	130.91	122.70
1	D	224	GLN	OE1-CD-NE2	-5.43	117.17	122.60
1	A	225	THR	N-CA-C	5.42	117.19	111.28
1	A	179	LEU	N-CA-C	5.40	118.09	111.82
1	A	70	HIS	CB-CG-CD2	-5.39	124.19	131.20
1	A	151	HIS	CA-CB-CG	-5.38	108.42	113.80
1	D	260	HIS	CB-CG-ND1	5.37	130.75	122.70
2	B	95	TRP	CG-CD2-CE3	5.37	139.27	133.90
2	E	50	GLU	CA-CB-CG	-5.34	103.42	114.10
1	D	151	HIS	CA-CB-CG	-5.32	108.48	113.80
1	D	255	GLN	OE1-CD-NE2	-5.31	117.29	122.60
1	A	173	GLU	CB-CG-CD	5.27	121.55	112.60
1	A	268	LYS	CG-CD-CE	-5.25	99.22	111.30
1	A	42	SER	N-CA-C	5.21	116.95	111.28
1	A	260	HIS	CB-CG-ND1	5.21	130.51	122.70
1	D	244	TRP	CE2-CD2-CG	-5.20	100.96	107.20
2	E	70	PHE	CA-CB-CG	5.20	119.00	113.80
2	B	50	GLU	CA-CB-CG	-5.20	103.70	114.10
1	A	208	PHE	CA-CB-CG	5.18	118.98	113.80
1	A	182	THR	CB-CA-C	5.17	117.58	110.34
1	D	72	GLN	OE1-CD-NE2	-5.16	117.44	122.60
1	A	220	ASP	CA-CB-CG	5.13	117.73	112.60
1	A	244	TRP	CE2-CD2-CG	-5.13	101.05	107.20
1	A	5	MET	CA-CB-CG	-5.12	103.86	114.10
2	B	95	TRP	CE2-CD2-CG	-5.12	101.06	107.20
1	D	39	ASP	CA-CB-CG	5.11	117.71	112.60
1	D	142	THR	CA-CB-OG1	-5.11	101.94	109.60
1	D	70	HIS	CB-CG-CD2	-5.10	124.57	131.20
2	E	13	HIS	CB-CG-ND1	5.09	130.34	122.70
1	D	42	SER	N-CA-C	5.09	116.83	111.28
1	D	274	TRP	CE2-CD2-CG	-5.07	101.11	107.20
1	A	39	ASP	CA-CB-CG	5.06	117.66	112.60
1	A	51	TRP	CE2-CD2-CG	-5.06	101.13	107.20
1	D	263	HIS	CB-CG-ND1	5.05	130.27	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	114	HIS	CB-CG-CD2	-5.04	124.65	131.20
2	E	95	TRP	CG-CD2-CE3	5.04	138.94	133.90
1	D	41	ALA	N-CA-C	5.03	116.76	111.28
1	D	51	TRP	CE2-CD2-CG	-5.02	101.17	107.20
2	E	95	TRP	CE2-CD2-CG	-5.02	101.17	107.20
1	A	266	LEU	CA-C-N	5.01	124.80	119.28
1	A	266	LEU	C-N-CA	5.01	124.80	119.28
1	A	192	HIS	CB-CG-CD2	-5.01	124.69	131.20
1	A	207	SER	N-CA-C	5.01	118.35	111.54

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	18	0
1	D	2247	0	2096	14	0
2	B	837	0	803	8	0
2	E	837	0	803	4	0
3	C	77	0	79	2	0
3	F	77	0	79	2	0
All	All	6322	0	5956	43	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 4.

All (43) 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:5:MET:HE1	1:D:7:TYR:CE2	2.32	0.65
2:E:36:GLU:HB3	2:E:83:ASN:HB2	1.79	0.65
2:B:36:GLU:HB3	2:B:83:ASN:HB2	1.79	0.64
1:A:5:MET:HE1	1:A:7:TYR:CE2	2.34	0.63
1:A:4:SER:HB2	1:A:6:ARG:HH12	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:4:SER:HB2	1:D:6:ARG:HH12	1.69	0.58
2:E:22:PHE:CE2	2:E:69:GLU:HG2	2.38	0.57
2:B:22:PHE:CE2	2:B:69:GLU:HG2	2.40	0.57
1:A:201:LEU:HD12	1:A:249:VAL:HG21	1.87	0.56
1:A:93:HIS:HD2	1:A:119:ASP:OD2	1.89	0.56
1:D:201:LEU:HD12	1:D:249:VAL:HG21	1.86	0.56
1:D:93:HIS:HD2	1:D:119:ASP:OD2	1.89	0.55
1:A:4:SER:HB2	1:A:6:ARG:NH1	2.25	0.52
1:D:210:PRO:O	1:D:263:HIS:HE1	1.91	0.52
2:E:25:CYS:HB2	2:E:39:LEU:HD21	1.92	0.52
1:A:210:PRO:O	1:A:263:HIS:HE1	1.93	0.52
2:B:25:CYS:HB2	2:B:39:LEU:HD21	1.93	0.51
1:D:4:SER:HB2	1:D:6:ARG:NH1	2.24	0.51
1:D:117:ALA:HB2	2:E:60:TRP:CE2	2.46	0.50
1:A:111:ARG:HD2	1:A:113:TYR:OH	2.10	0.49
1:D:111:ARG:HD2	1:D:113:TYR:OH	2.12	0.49
1:A:232:GLU:OE2	2:B:6:LYS:NZ	2.42	0.49
1:D:165:VAL:O	1:D:169:ARG:HG2	2.12	0.48
1:A:49:ALA:O	1:A:52:ILE:HG22	2.14	0.48
1:A:182:THR:HG23	1:A:265:GLY:HA2	1.97	0.47
1:A:169:ARG:HA	1:A:172:LEU:HD12	1.96	0.47
1:D:49:ALA:O	1:D:52:ILE:HG22	2.14	0.47
1:D:21:ARG:HH21	1:D:23:ILE:HD11	1.80	0.46
1:D:146:LYS:NZ	3:F:9:VAL:O	2.49	0.46
1:A:6:ARG:NH2	1:A:113:TYR:CE1	2.83	0.46
1:D:6:ARG:NH2	1:D:113:TYR:CE1	2.84	0.46
2:B:73:THR:O	2:B:97:ARG:NH2	2.49	0.46
1:D:11:SER:HA	1:D:21:ARG:O	2.15	0.46
1:A:11:SER:HA	1:A:21:ARG:O	2.16	0.45
3:F:3:PHE:CZ	3:F:5:TYR:HB2	2.53	0.44
1:A:21:ARG:HH21	1:A:23:ILE:HD11	1.82	0.44
3:C:3:PHE:CZ	3:C:5:TYR:HB2	2.53	0.43
2:B:81:ARG:HG3	2:B:92:ILE:HG12	2.00	0.42
1:A:84:TYR:HB3	1:A:139:ALA:HB1	2.02	0.42
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.56	0.41
1:A:146:LYS:NZ	3:C:9:VAL:O	2.53	0.41
1:A:203:CYS:HB2	1:A:217:TRP:CZ2	2.56	0.40
2:B:37:VAL:HG22	2:B:82:VAL:HG22	2.03	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%)	267 (98%)	6 (2%)	0	100	100
1	D	273/275 (99%)	269 (98%)	4 (2%)	0	100	100
2	B	98/100 (98%)	97 (99%)	1 (1%)	0	100	100
2	E	98/100 (98%)	97 (99%)	1 (1%)	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%)	744 (98%)	12 (2%)	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%)	226 (98%)	5 (2%)	45	73
1	D	231/231 (100%)	224 (97%)	7 (3%)	36	64
2	B	95/95 (100%)	93 (98%)	2 (2%)	47	74
2	E	95/95 (100%)	93 (98%)	2 (2%)	47	74
3	C	8/8 (100%)	8 (100%)	0	100	100
3	F	8/8 (100%)	8 (100%)	0	100	100
All	All	668/668 (100%)	652 (98%)	16 (2%)	43	70

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	MET
1	A	11	SER
1	A	58	GLU
1	A	75	ARG
1	A	182	THR
2	B	9	VAL
2	B	34	ASP
1	D	5	MET
1	D	11	SER
1	D	58	GLU
1	D	75	ARG
1	D	105	SER
1	D	110	LEU
1	D	169	ARG
2	E	9	VAL
2	E	34	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	HIS
1	A	180	GLN
1	A	263	HIS
1	D	93	HIS
1	D	263	HIS
2	E	83	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

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.