



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:25 AM UTC

PDB ID : 2HLA / pdb_00002hla
Title : SPECIFICITY POCKETS FOR THE SIDE CHAINS OF PEPTIDE ANTI-GENS IN HLA-AW68
Authors : Garrett, T.P.J.; Saper, M.A.; Wiley, D.C.
Deposited on : 1989-10-05
Resolution : 2.60 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

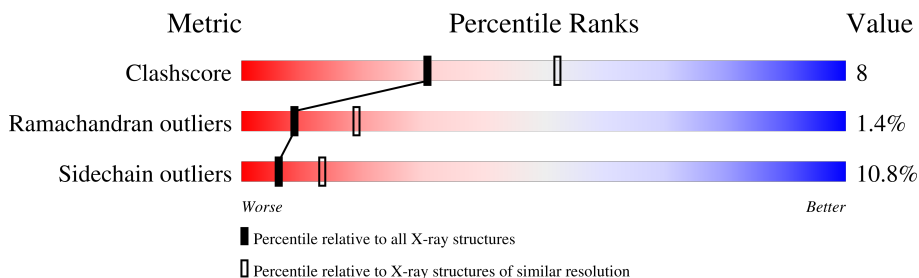
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)

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	270	
2	B	99	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3033 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-Aw68).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	270	Total	C	N	O	S	0	0	0
			2194	1362	399	423	10			

- 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	0	0
			829	528	140	158	3			

- Molecule 3 is water.

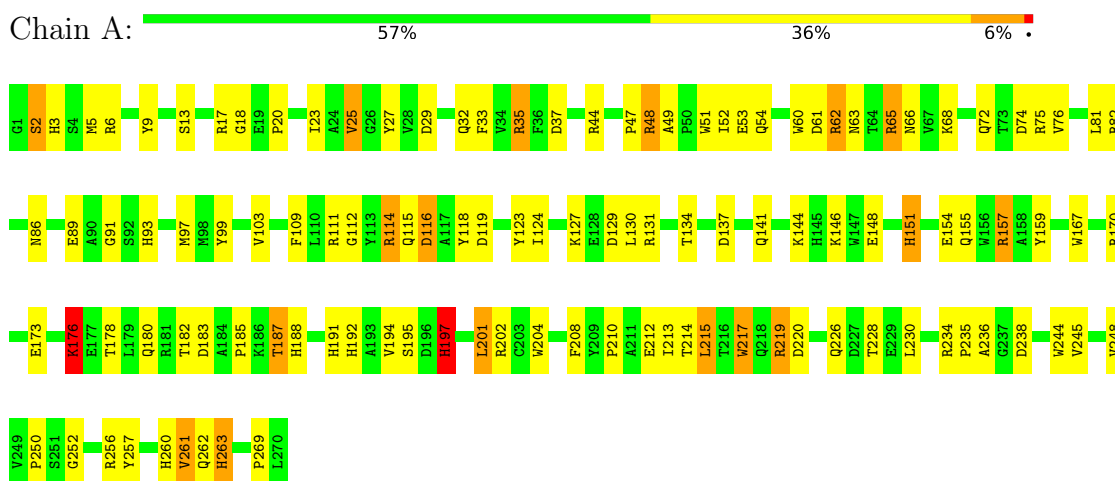
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	6	Total	O	0	0
			6	6		
3	B	4	Total	O	0	0
			4	4		

3 Residue-property plots

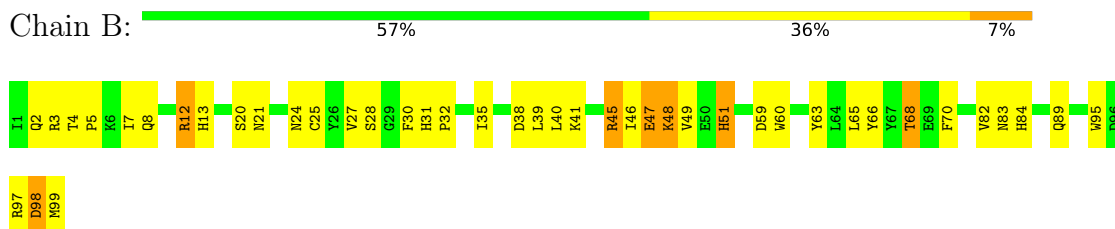
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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-Aw68)



• Molecule 2: BETA 2-MICROGLOBULIN



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	60.00Å 80.66Å 113.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.60	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-2.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.173 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3033	wwPDB-VP
Average B, all atoms (Å ²)	20.0	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	1.35	14/2253 (0.6%)	2.17	84/3056 (2.7%)
2	B	1.38	8/852 (0.9%)	2.13	42/1152 (3.6%)
All	All	1.36	22/3105 (0.7%)	2.16	126/4208 (3.0%)

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	A	0	1

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	263	HIS	CD2-NE2	-7.55	1.29	1.37
1	A	76	VAL	CA-CB	7.44	1.63	1.54
2	B	51	HIS	CD2-NE2	-7.13	1.30	1.37
1	A	260	HIS	CD2-NE2	-6.98	1.30	1.37
2	B	31	HIS	CD2-NE2	-6.59	1.30	1.37
2	B	84	HIS	CD2-NE2	-6.54	1.30	1.37
2	B	13	HIS	CD2-NE2	-6.52	1.30	1.37
1	A	192	HIS	CD2-NE2	-6.46	1.30	1.37
1	A	3	HIS	CD2-NE2	-6.35	1.30	1.37
1	A	188	HIS	CD2-NE2	-6.26	1.30	1.37
1	A	261	VAL	CA-CB	5.84	1.61	1.54
2	B	51	HIS	CG-ND1	-5.83	1.31	1.38
1	A	187	THR	CA-CB	5.77	1.62	1.53
2	B	13	HIS	CG-ND1	-5.68	1.32	1.38
1	A	191	HIS	CD2-NE2	-5.47	1.31	1.37
2	B	31	HIS	CG-ND1	-5.45	1.32	1.38
1	A	3	HIS	CG-ND1	-5.22	1.32	1.38
1	A	197	HIS	CD2-NE2	-5.19	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	93	HIS	CD2-NE2	-5.19	1.32	1.37
1	A	188	HIS	CG-ND1	-5.09	1.32	1.38
2	B	41	LYS	CA-CB	-5.08	1.46	1.53
1	A	151	HIS	CD2-NE2	-5.05	1.32	1.37

All (126) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	74	ASP	CA-CB-CG	19.54	132.14	112.60
1	A	197	HIS	CA-CB-CG	12.97	126.77	113.80
1	A	114	ARG	NE-CZ-NH1	-10.09	111.41	121.50
1	A	114	ARG	NE-CZ-NH2	9.73	127.96	119.20
1	A	151	HIS	CA-CB-CG	-9.54	104.26	113.80
2	B	45	ARG	NE-CZ-NH1	-9.06	112.44	121.50
1	A	62	ARG	NE-CZ-NH1	-9.05	112.45	121.50
1	A	63	ASN	OD1-CG-ND2	-8.89	113.71	122.60
2	B	24	ASN	OD1-CG-ND2	-8.80	113.80	122.60
1	A	235	PRO	CA-C-O	-8.57	112.06	121.23
1	A	62	ARG	NE-CZ-NH2	8.56	126.90	119.20
1	A	116	ASP	CA-CB-CG	8.54	121.14	112.60
2	B	84	HIS	ND1-CG-CD2	8.46	114.56	106.10
2	B	45	ARG	NE-CZ-NH2	8.45	126.80	119.20
1	A	170	ARG	NE-CZ-NH1	-7.76	113.74	121.50
2	B	21	ASN	CA-CB-CG	7.66	120.26	112.60
1	A	66	ASN	OD1-CG-ND2	-7.58	115.02	122.60
1	A	137	ASP	CA-CB-CG	7.56	120.16	112.60
1	A	72	GLN	OE1-CD-NE2	-7.42	115.18	122.60
1	A	23	ILE	CB-CA-C	-7.40	100.43	110.42
1	A	252	GLY	N-CA-C	-7.38	105.08	114.37
1	A	141	GLN	OE1-CD-NE2	-7.20	115.41	122.60
1	A	63	ASN	CB-CG-ND2	7.19	127.19	116.40
1	A	131	ARG	NE-CZ-NH1	-7.08	114.42	121.50
1	A	61	ASP	CA-CB-CG	-6.96	105.64	112.60
1	A	226	GLN	OE1-CD-NE2	-6.95	115.65	122.60
1	A	256	ARG	NE-CZ-NH1	-6.93	114.57	121.50
1	A	170	ARG	CG-CD-NE	-6.88	96.86	112.00
1	A	219	ARG	NE-CZ-NH1	-6.85	114.65	121.50
2	B	12	ARG	CB-CG-CD	-6.78	95.70	111.30
1	A	68	LYS	CA-CB-CG	-6.75	100.60	114.10
2	B	97	ARG	CB-CG-CD	-6.71	95.87	111.30
1	A	48	ARG	NE-CZ-NH1	-6.71	114.80	121.50
2	B	59	ASP	CA-CB-CG	6.69	119.29	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	18	GLY	N-CA-C	6.66	120.22	111.70
1	A	204	TRP	CE2-CD2-CE3	6.61	125.41	118.80
1	A	65	ARG	CB-CA-C	-6.59	98.44	110.63
2	B	84	HIS	CG-CD2-NE2	-6.57	100.63	107.20
2	B	38	ASP	CA-CB-CG	6.53	119.13	112.60
1	A	111	ARG	CA-CB-CG	6.42	126.94	114.10
1	A	167	TRP	CG-CD2-CE3	6.42	140.32	133.90
1	A	236	ALA	N-CA-C	-6.28	104.51	111.36
2	B	60	TRP	CG-CD2-CE3	6.23	140.13	133.90
1	A	65	ARG	NE-CZ-NH2	6.21	124.79	119.20
1	A	155	GLN	OE1-CD-NE2	-6.21	116.39	122.60
2	B	41	LYS	CB-CG-CD	-6.19	97.07	111.30
1	A	29	ASP	CB-CA-C	-6.18	109.46	116.63
2	B	89	GLN	CA-C-N	6.15	126.17	119.90
2	B	89	GLN	C-N-CA	6.15	126.17	119.90
1	A	129	ASP	CA-CB-CG	6.11	118.70	112.60
2	B	31	HIS	ND1-CG-CD2	6.07	112.17	106.10
2	B	98	ASP	CA-CB-CG	6.06	118.66	112.60
2	B	83	ASN	OD1-CG-ND2	-6.04	116.56	122.60
1	A	127	LYS	CB-CG-CD	-6.03	97.43	111.30
1	A	195	SER	CA-C-N	-5.98	113.76	122.19
1	A	195	SER	C-N-CA	-5.98	113.76	122.19
1	A	202	ARG	NE-CZ-NH1	-5.97	115.53	121.50
1	A	9	TYR	CA-CB-CG	5.97	124.64	113.90
1	A	33	PHE	CA-CB-CG	5.94	119.74	113.80
1	A	65	ARG	NE-CZ-NH1	-5.94	115.56	121.50
1	A	6	ARG	NE-CZ-NH1	-5.92	115.58	121.50
1	A	248	VAL	N-CA-C	-5.89	101.20	109.21
2	B	21	ASN	CB-CG-ND2	-5.88	107.57	116.40
1	A	35	ARG	NE-CZ-NH1	-5.86	115.64	121.50
2	B	13	HIS	ND1-CG-CD2	5.84	111.94	106.10
1	A	191	HIS	ND1-CG-CD2	5.77	111.87	106.10
1	A	5	MET	CA-CB-CG	-5.75	102.60	114.10
2	B	31	HIS	O-C-N	-5.73	114.73	121.32
2	B	24	ASN	CB-CG-ND2	5.72	124.98	116.40
1	A	183	ASP	CA-CB-CG	5.71	118.31	112.60
1	A	202	ARG	NE-CZ-NH2	5.69	124.32	119.20
1	A	157	ARG	N-CA-C	-5.68	104.29	111.11
1	A	119	ASP	N-CA-C	5.67	119.01	111.30
1	A	191	HIS	CB-CG-CD2	-5.63	123.88	131.20
2	B	47	GLU	N-CA-C	-5.60	102.18	110.46
2	B	60	TRP	CE2-CD2-CG	-5.58	100.51	107.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	244	TRP	CB-CG-CD1	-5.54	118.60	126.90
2	B	49	VAL	CA-C-O	-5.51	114.52	120.36
1	A	244	TRP	CG-CD2-CE3	5.51	139.41	133.90
1	A	44	ARG	NE-CZ-NH1	-5.50	116.00	121.50
1	A	75	ARG	NE-CZ-NH1	-5.50	116.00	121.50
2	B	95	TRP	CE2-CD2-CG	-5.49	100.61	107.20
2	B	24	ASN	CA-CB-CG	5.48	118.08	112.60
1	A	112	GLY	CA-C-O	-5.47	115.78	121.63
1	A	215	LEU	N-CA-CB	-5.39	102.11	110.57
1	A	134	THR	CA-C-O	5.39	126.84	120.70
1	A	66	ASN	CA-C-O	-5.36	114.81	120.55
1	A	220	ASP	CA-CB-CG	5.36	117.96	112.60
1	A	60	TRP	CG-CD2-CE3	5.35	139.25	133.90
1	A	262	GLN	OE1-CD-NE2	-5.35	117.25	122.60
2	B	97	ARG	N-CA-CB	-5.34	101.70	110.14
2	B	40	LEU	N-CA-C	5.34	117.43	108.73
2	B	84	HIS	CB-CG-CD2	-5.33	124.28	131.20
1	A	86	ASN	CA-CB-CG	-5.31	107.29	112.60
2	B	66	TYR	CB-CG-CD1	-5.31	112.84	120.80
2	B	48	LYS	CA-CB-CG	5.30	124.69	114.10
2	B	70	PHE	CA-CB-CG	5.25	119.05	113.80
1	A	115	GLN	OE1-CD-NE2	-5.24	117.36	122.60
1	A	141	GLN	N-CA-C	-5.22	105.67	111.36
1	A	167	TRP	CG-CD1-NE1	-5.22	103.42	110.20
1	A	2	SER	CA-C-O	-5.21	115.77	121.14
1	A	37	ASP	O-C-N	-5.21	117.00	123.36
2	B	49	VAL	CB-CA-C	-5.21	102.72	110.33
1	A	260	HIS	CB-CG-CD2	-5.19	124.46	131.20
1	A	146	LYS	N-CA-C	-5.18	105.64	111.28
1	A	217	TRP	CG-CD2-CE3	5.17	139.07	133.90
1	A	109	PHE	CA-CB-CG	5.17	118.97	113.80
1	A	188	HIS	ND1-CG-CD2	5.17	111.27	106.10
2	B	20	SER	N-CA-C	-5.17	102.83	110.48
1	A	167	TRP	CE2-CD2-CG	-5.16	101.01	107.20
1	A	25	VAL	N-CA-C	-5.14	101.22	108.27
2	B	68	THR	CA-CB-OG1	-5.14	101.89	109.60
1	A	35	ARG	O-C-N	5.14	129.14	123.33
1	A	17	ARG	N-CA-C	-5.13	99.87	110.80
1	A	236	ALA	CA-C-O	5.12	125.85	120.42
2	B	51	HIS	ND1-CG-CD2	5.12	111.22	106.10
1	A	25	VAL	CG1-CB-CG2	-5.10	99.57	110.80
2	B	97	ARG	N-CA-C	5.10	118.28	111.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	47	PRO	CA-C-O	5.09	127.42	121.31
1	A	228	THR	O-C-N	-5.08	116.76	123.16
2	B	30	PHE	CA-CB-CG	5.08	118.88	113.80
2	B	99	MET	CG-SD-CE	-5.08	89.73	100.90
1	A	263	HIS	CB-CG-CD2	-5.03	124.67	131.20
2	B	3	ARG	CA-C-N	5.02	127.62	120.49
2	B	3	ARG	C-N-CA	5.02	127.62	120.49
2	B	31	HIS	CB-CG-CD2	-5.01	124.68	131.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	159	TYR	Sidechain

5.2 Too-close contacts [i](#)

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	2194	0	2045	35	0
2	B	829	0	794	16	0
3	A	6	0	0	0	0
3	B	4	0	0	0	0
All	All	3033	0	2839	48	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 8.

All (48) 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:62:ARG:HD2	1:A:65:ARG:NH1	2.07	0.69
1:A:13:SER:HA	1:A:20:PRO:HB3	1.76	0.66
2:B:2:GLN:HG2	2:B:32:PRO:HD3	1.78	0.66
1:A:62:ARG:HD2	1:A:65:ARG:HH11	1.60	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:7:ILE:HG12	2:B:82:VAL:HG21	1.82	0.61
2:B:25:CYS:HB2	2:B:39:LEU:HD21	1.83	0.59
1:A:173:GLU:O	1:A:176:LYS:HD3	2.03	0.59
1:A:238:ASP:HB3	2:B:12:ARG:HH11	1.67	0.58
2:B:46:ILE:HD13	2:B:68:THR:HG21	1.86	0.57
2:B:47:GLU:O	2:B:48:LYS:HB3	2.06	0.54
1:A:2:SER:HB2	1:A:103:VAL:O	2.07	0.54
1:A:130:LEU:HB3	1:A:157:ARG:HG3	1.90	0.54
1:A:187:THR:HG21	1:A:261:VAL:HG21	1.90	0.54
1:A:25:VAL:HG12	1:A:27:TYR:CE2	2.43	0.54
1:A:185:PRO:HB3	1:A:208:PHE:HB3	1.90	0.53
2:B:45:ARG:HH12	2:B:47:GLU:CD	2.17	0.52
1:A:238:ASP:HB3	2:B:12:ARG:HE	1.74	0.51
1:A:62:ARG:HA	1:A:65:ARG:NH1	2.26	0.51
1:A:25:VAL:HG12	1:A:27:TYR:HE2	1.76	0.50
1:A:49:ALA:O	1:A:52:ILE:HG22	2.12	0.50
1:A:32:GLN:NE2	1:A:48:ARG:HG3	2.27	0.49
1:A:82:ARG:NH1	1:A:91:GLY:O	2.46	0.49
1:A:176:LYS:HA	1:A:180:GLN:HG3	1.95	0.49
2:B:27:VAL:HG11	2:B:35:ILE:HD13	1.94	0.48
1:A:213:ILE:HD12	1:A:263:HIS:HB2	1.97	0.47
1:A:230:LEU:HD12	1:A:245:VAL:HB	1.97	0.46
1:A:219:ARG:HG3	1:A:257:TYR:CZ	2.51	0.46
1:A:116:ASP:OD2	1:A:123:TYR:HB3	2.15	0.46
1:A:51:TRP:CD1	1:A:51:TRP:H	2.33	0.46
1:A:144:LYS:O	1:A:148:GLU:HG3	2.16	0.46
1:A:51:TRP:O	1:A:54:GLN:HG2	2.15	0.45
1:A:114:ARG:NH1	1:A:124:ILE:HG23	2.31	0.45
1:A:123:TYR:HD1	1:A:124:ILE:HG22	1.82	0.45
2:B:25:CYS:HB2	2:B:39:LEU:CD2	2.47	0.45
2:B:51:HIS:HA	2:B:65:LEU:O	2.17	0.45
1:A:215:LEU:CD1	1:A:261:VAL:HG22	2.48	0.44
1:A:210:PRO:O	1:A:263:HIS:HE1	2.00	0.44
1:A:234:ARG:HD3	2:B:8:GLN:OE1	2.19	0.43
1:A:201:LEU:HB3	1:A:217:TRP:CH2	2.53	0.43
2:B:48:LYS:HB3	2:B:48:LYS:HE3	1.79	0.42
1:A:81:LEU:HD23	1:A:118:TYR:CD1	2.55	0.42
1:A:82:ARG:NE	1:A:89:GLU:HA	2.34	0.42
1:A:215:LEU:HD12	1:A:261:VAL:HG22	2.00	0.42
2:B:28:SER:HA	2:B:63:TYR:HA	2.01	0.42
1:A:197:HIS:O	1:A:250:PRO:HA	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:MET:HE3	1:A:99:TYR:CD1	2.57	0.40
2:B:27:VAL:HG11	2:B:35:ILE:CD1	2.51	0.40
2:B:4:THR:HA	2:B:5:PRO:HD2	1.97	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	254/270 (94%)	232 (91%)	17 (7%)	5 (2%)	6	12
2	B	97/99 (98%)	90 (93%)	7 (7%)	0	100	100
All	All	351/369 (95%)	322 (92%)	24 (7%)	5 (1%)	9	19

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	269	PRO
1	A	194	VAL
1	A	197	HIS
1	A	176	LYS
1	A	151	HIS

5.3.2 Protein sidechains [i](#)

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	89/227 (39%)	80 (90%)	9 (10%)	7	15
2	B	4/94 (4%)	3 (75%)	1 (25%)	0	1
All	All	93/321 (29%)	83 (89%)	10 (11%)	6	13

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	ARG
1	A	53	GLU
1	A	154	GLU
1	A	176	LYS
1	A	178	THR
1	A	182	THR
1	A	201	LEU
1	A	212	GLU
1	A	214	THR
2	B	98	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	141	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

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.