



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:11 PM UTC

PDB ID : 3HLA / pdb_00003hla
Title : HUMAN CLASS I HISTOCOMPATIBILITY ANTIGEN A2.1
Authors : Saper, M.A.; Bjorkman, P.J.; Wiley, D.C.
Deposited on : 1989-10-06
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

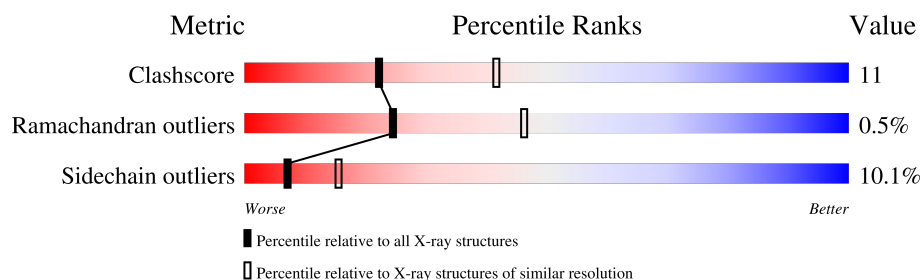
1 Overall quality at a glance

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 3064 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-A2.1) (ALPHA CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	270	Total	C	N	O	S	0	0	0
			2189	1363	399	418	9			

- 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.

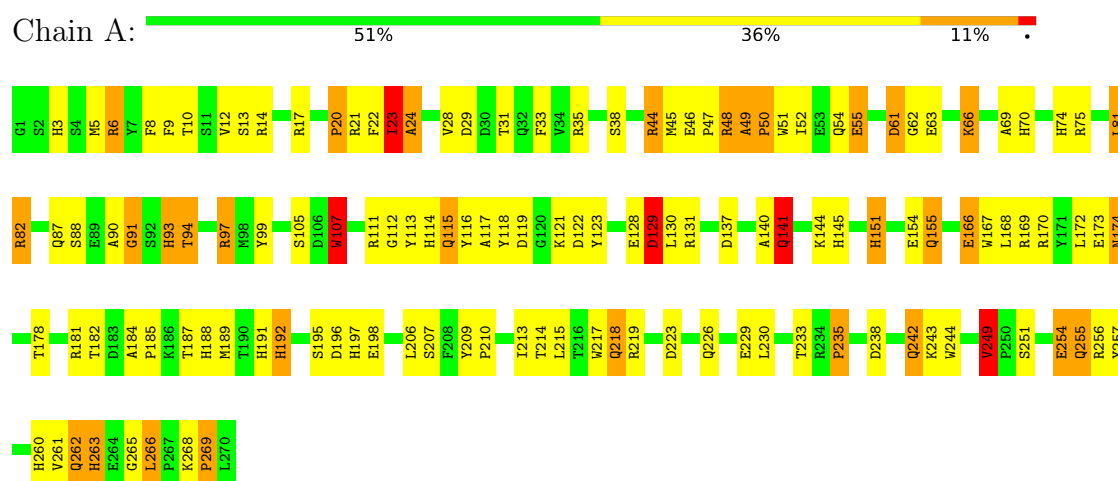
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	30	Total	O	0	0
			30	30		
3	B	16	Total	O	0	0
			16	16		

3 Residue-property plots

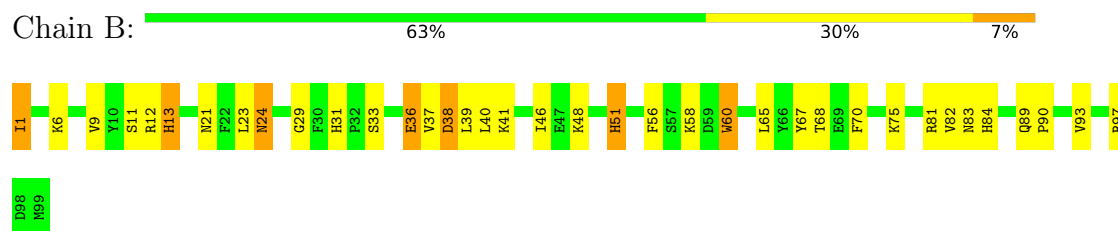
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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-A2.1) (ALPHA CHAIN)



• 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 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.35Å 80.40Å 56.49Å 90.00° 120.42° 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}	0.169 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3064	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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	1.39	19/2251 (0.8%)	2.14	94/3054 (3.1%)
2	B	1.32	8/852 (0.9%)	2.01	14/1152 (1.2%)
All	All	1.37	27/3103 (0.9%)	2.11	108/4206 (2.6%)

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	70	HIS	CD2-NE2	-7.57	1.29	1.37
2	B	13	HIS	CD2-NE2	-7.15	1.29	1.37
1	A	114	HIS	CD2-NE2	-7.10	1.30	1.37
1	A	93	HIS	CD2-NE2	-6.98	1.30	1.37
2	B	84	HIS	CD2-NE2	-6.80	1.30	1.37
1	A	145	HIS	CD2-NE2	-6.48	1.30	1.37
2	B	84	HIS	CG-ND1	-6.39	1.31	1.38
1	A	70	HIS	CG-ND1	-6.30	1.31	1.38
1	A	263	HIS	CD2-NE2	-6.29	1.30	1.37
1	A	191	HIS	CD2-NE2	-6.20	1.31	1.37
2	B	31	HIS	CG-ND1	-6.20	1.31	1.38
1	A	74	HIS	CD2-NE2	-5.97	1.31	1.37
1	A	192	HIS	CD2-NE2	-5.94	1.31	1.37
1	A	260	HIS	CD2-NE2	-5.84	1.31	1.37
1	A	188	HIS	CD2-NE2	-5.81	1.31	1.37
1	A	93	HIS	CG-ND1	-5.75	1.31	1.38
1	A	197	HIS	CD2-NE2	-5.70	1.31	1.37
2	B	13	HIS	CG-ND1	-5.62	1.32	1.38
1	A	3	HIS	CD2-NE2	-5.59	1.31	1.37
1	A	91	GLY	CA-C	5.54	1.56	1.52
1	A	151	HIS	CD2-NE2	-5.50	1.31	1.37
1	A	151	HIS	CG-ND1	-5.39	1.32	1.38
2	B	51	HIS	CD2-NE2	-5.32	1.31	1.37
1	A	3	HIS	CG-ND1	-5.16	1.32	1.38
2	B	48	LYS	CA-CB	5.15	1.61	1.53
2	B	31	HIS	CD2-NE2	-5.08	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	48	ARG	N-CA	-5.02	1.40	1.46

All (108) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	38	ASP	CA-CB-CG	16.72	129.32	112.60
1	A	249	VAL	N-CA-CB	-13.83	99.92	111.67
2	B	24	ASN	CB-CG-ND2	9.64	130.86	116.40
2	B	24	ASN	OD1-CG-ND2	-9.39	113.21	122.60
1	A	93	HIS	CA-CB-CG	-9.18	104.62	113.80
1	A	260	HIS	CA-CB-CG	-8.76	105.04	113.80
1	A	137	ASP	CA-CB-CG	8.59	121.19	112.60
1	A	269	PRO	N-CA-CB	8.58	112.26	103.25
2	B	56	PHE	CA-CB-CG	8.53	122.33	113.80
2	B	89	GLN	N-CA-CB	-8.45	102.23	111.10
1	A	145	HIS	CA-CB-CG	-8.41	105.39	113.80
1	A	242	GLN	OE1-CD-NE2	-8.38	114.22	122.60
1	A	233	THR	CA-CB-OG1	-8.03	97.55	109.60
1	A	55	GLU	CB-CG-CD	7.57	125.47	112.60
1	A	174	ASN	OD1-CG-ND2	-7.56	115.04	122.60
1	A	28	VAL	N-CA-C	-7.46	96.74	107.77
1	A	38	SER	CA-CB-OG	-7.44	96.23	111.10
1	A	196	ASP	CA-CB-CG	7.42	120.02	112.60
1	A	46	GLU	CA-C-N	7.14	127.06	119.85
1	A	46	GLU	C-N-CA	7.14	127.06	119.85
1	A	97	ARG	NE-CZ-NH2	6.82	125.34	119.20
1	A	119	ASP	N-CA-C	6.73	120.79	111.90
1	A	218	GLN	OE1-CD-NE2	-6.72	115.88	122.60
1	A	33	PHE	CA-CB-CG	6.67	120.47	113.80
1	A	105	SER	CA-CB-OG	-6.67	97.76	111.10
1	A	31	THR	CA-CB-OG1	-6.59	99.71	109.60
2	B	89	GLN	CA-CB-CG	6.58	127.26	114.10
1	A	29	ASP	CA-CB-CG	6.47	119.07	112.60
1	A	49	ALA	CA-C-N	6.35	127.77	119.84
1	A	49	ALA	C-N-CA	6.35	127.77	119.84
1	A	155	GLN	OE1-CD-NE2	-6.34	116.25	122.60
1	A	219	ARG	NE-CZ-NH1	-6.34	115.16	121.50
1	A	75	ARG	N-CA-C	-6.26	104.38	111.14
1	A	75	ARG	NE-CZ-NH2	6.21	124.79	119.20
1	A	61	ASP	CA-CB-CG	-6.16	106.44	112.60
1	A	155	GLN	N-CA-C	-6.15	104.49	111.07
2	B	84	HIS	ND1-CG-CD2	6.14	112.24	106.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	181	ARG	O-C-N	6.13	131.11	123.14
1	A	188	HIS	CA-CB-CG	6.11	119.91	113.80
1	A	94	THR	O-C-N	-6.08	116.09	123.27
1	A	111	ARG	CA-CB-CG	6.05	126.20	114.10
1	A	47	PRO	N-CA-CB	5.99	108.90	103.33
1	A	141	GLN	OE1-CD-NE2	-5.98	116.62	122.60
1	A	8	PHE	CA-CB-CG	-5.96	107.84	113.80
1	A	21	ARG	NE-CZ-NH2	5.95	124.56	119.20
2	B	60	TRP	CG-CD1-NE1	-5.94	102.48	110.20
1	A	262	GLN	CA-CB-CG	5.93	125.96	114.10
1	A	151	HIS	CA-CB-CG	-5.90	107.90	113.80
1	A	105	SER	N-CA-C	5.89	118.48	111.71
1	A	122	ASP	CA-CB-CG	5.86	118.46	112.60
1	A	244	TRP	CG-CD2-CE3	5.84	139.74	133.90
1	A	47	PRO	CA-N-CD	-5.83	103.84	112.00
1	A	235	PRO	N-CA-CB	-5.79	98.55	103.36
1	A	129	ASP	N-CA-CB	-5.77	101.40	110.46
1	A	69	ALA	O-C-N	5.76	128.08	122.09
2	B	24	ASN	CA-CB-CG	5.73	118.33	112.60
2	B	97	ARG	CB-CG-CD	-5.73	98.12	111.30
1	A	17	ARG	NE-CZ-NH2	5.72	124.35	119.20
1	A	23	ILE	CB-CA-C	-5.64	102.21	110.62
1	A	107	TRP	CB-CG-CD1	-5.63	118.45	126.90
1	A	235	PRO	CB-CA-C	5.62	117.68	111.11
1	A	23	ILE	CA-CB-CG1	5.59	119.91	110.40
1	A	90	ALA	N-CA-C	5.58	119.28	111.30
1	A	88	SER	O-C-N	5.55	129.12	122.79
2	B	60	TRP	CD1-CG-CD2	5.54	115.17	106.30
1	A	255	GLN	OE1-CD-NE2	-5.53	117.07	122.60
1	A	20	PRO	CB-CA-C	5.50	118.50	110.63
1	A	254	GLU	CA-CB-CG	5.50	125.09	114.10
1	A	144	LYS	CB-CG-CD	-5.49	98.67	111.30
1	A	35	ARG	NE-CZ-NH1	-5.45	116.06	121.50
1	A	137	ASP	N-CA-CB	-5.43	102.15	111.21
1	A	29	ASP	CB-CA-C	-5.41	110.32	116.54
1	A	6	ARG	NE-CZ-NH1	-5.39	116.11	121.50
1	A	44	ARG	CA-CB-CG	5.37	124.85	114.10
1	A	189	MET	CB-CG-SD	-5.32	96.74	112.70
1	A	244	TRP	CB-CG-CD1	-5.31	118.93	126.90
1	A	24	ALA	N-CA-C	-5.31	99.79	108.76
1	A	262	GLN	OE1-CD-NE2	-5.30	117.30	122.60
1	A	82	ARG	CB-CG-CD	-5.26	99.20	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	60	TRP	CE2-CD2-CG	-5.24	100.92	107.20
1	A	75	ARG	NE-CZ-NH1	-5.21	116.30	121.50
1	A	112	GLY	O-C-N	-5.20	118.73	123.36
1	A	256	ARG	NE-CZ-NH2	5.20	123.88	119.20
1	A	50	PRO	CA-N-CD	-5.19	104.73	112.00
1	A	238	ASP	CA-CB-CG	5.19	117.79	112.60
1	A	217	TRP	CG-CD1-NE1	-5.19	103.45	110.20
2	B	58	LYS	N-CA-C	5.19	117.68	111.71
1	A	192	HIS	O-C-N	-5.18	117.21	123.27
1	A	244	TRP	CE2-CD2-CG	-5.17	101.00	107.20
1	A	256	ARG	NE-CZ-NH1	-5.17	116.33	121.50
1	A	244	TRP	CG-CD1-NE1	-5.16	103.49	110.20
1	A	82	ARG	NE-CZ-NH2	5.16	123.84	119.20
1	A	196	ASP	N-CA-CB	-5.15	102.29	110.28
1	A	17	ARG	O-C-N	5.13	129.42	123.41
2	B	11	SER	CB-CA-C	-5.13	100.86	109.48
1	A	209	TYR	O-C-N	-5.13	116.11	122.16
1	A	75	ARG	CA-CB-CG	5.12	124.34	114.10
1	A	10	THR	CA-CB-OG1	-5.11	101.93	109.60
1	A	233	THR	CA-CB-CG2	5.08	119.13	110.50
1	A	107	TRP	CA-CB-CG	5.08	123.24	113.60
1	A	137	ASP	O-C-N	-5.07	116.98	122.92
1	A	90	ALA	CA-C-O	5.07	127.32	121.34
1	A	130	LEU	N-CA-CB	-5.07	104.56	112.41
1	A	51	TRP	CE2-CD2-CG	-5.06	101.13	107.20
1	A	191	HIS	CA-CB-CG	5.04	118.83	113.80
1	A	22	PHE	CA-C-N	5.03	130.00	123.06
1	A	22	PHE	C-N-CA	5.03	130.00	123.06
1	A	66	LYS	CB-CG-CD	-5.02	99.76	111.30

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	2189	0	2023	46	0
2	B	829	0	794	18	0
3	A	30	0	0	2	0
3	B	16	0	0	0	0
All	All	3064	0	2817	62	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 11.

All (62) 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:214:THR:HB	1:A:262:GLN:HB2	1.67	0.76
2:B:24:ASN:HB3	2:B:65:LEU:HD11	1.68	0.74
1:A:187:THR:HG21	1:A:261:VAL:HG21	1.74	0.70
1:A:218:GLN:HG2	1:A:223:ASP:HA	1.75	0.68
1:A:170:ARG:HH11	1:A:174:ASN:HD21	1.44	0.66
1:A:6:ARG:HD3	1:A:113:TYR:OH	1.97	0.64
1:A:49:ALA:O	1:A:52:ILE:HG22	2.01	0.60
2:B:36:GLU:HG3	2:B:83:ASN:HB3	1.87	0.56
2:B:83:ASN:HB2	2:B:90:PRO:HB3	1.86	0.56
1:A:81:LEU:HD23	1:A:118:TYR:CD1	2.41	0.56
1:A:45:MET:HG2	1:A:63:GLU:HB3	1.88	0.55
1:A:151:HIS:O	1:A:155:GLN:HG2	2.06	0.55
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.88	0.55
1:A:9:PHE:HE1	1:A:99:TYR:CE1	2.25	0.54
1:A:249:VAL:HG22	1:A:257:TYR:CE2	2.43	0.54
1:A:63:GLU:OE2	1:A:66:LYS:NZ	2.43	0.52
1:A:230:LEU:HD11	1:A:243:LYS:HE3	1.92	0.52
1:A:107:TRP:CZ2	1:A:172:LEU:HD13	2.44	0.52
2:B:6:LYS:HE3	2:B:29:GLY:HA3	1.90	0.52
2:B:38:ASP:OD2	2:B:81:ARG:NH2	2.43	0.52
1:A:129:ASP:OD1	1:A:131:ARG:HD2	2.10	0.51
1:A:44:ARG:HH12	1:A:61:ASP:CG	2.18	0.51
1:A:192:HIS:HB3	3:A:905:HOH:O	2.10	0.51
1:A:82:ARG:HG3	1:A:87:GLN:HB2	1.93	0.51
1:A:167:TRP:CZ3	1:A:170:ARG:HD3	2.46	0.50
1:A:82:ARG:NH1	1:A:91:GLY:O	2.40	0.50
1:A:184:ALA:HB2	1:A:265:GLY:O	2.12	0.50
1:A:206:LEU:HD23	1:A:242:GLN:HB3	1.94	0.49
1:A:185:PRO:HD2	1:A:266:LEU:HD13	1.93	0.49
1:A:12:VAL:HG22	1:A:94:THR:HG23	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:SER:HA	1:A:20:PRO:HB3	1.93	0.48
1:A:6:ARG:NH2	1:A:113:TYR:OH	2.45	0.48
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.48	0.47
1:A:182:THR:HG22	1:A:210:PRO:HD3	1.97	0.46
2:B:21:ASN:HB3	2:B:70:PHE:CE1	2.51	0.46
1:A:14:ARG:NH1	3:A:926:HOH:O	2.47	0.45
1:A:121:LYS:HE3	1:A:121:LYS:HB2	1.71	0.44
2:B:41:LYS:HB2	2:B:46:ILE:HD11	2.00	0.44
1:A:62:GLY:O	1:A:66:LYS:HG3	2.18	0.43
2:B:23:LEU:O	2:B:67:TYR:HA	2.18	0.43
1:A:93:HIS:HB3	1:A:118:TYR:CE1	2.54	0.43
1:A:23:ILE:HG22	1:A:24:ALA:N	2.34	0.42
1:A:206:LEU:CD2	1:A:242:GLN:HB3	2.49	0.42
2:B:12:ARG:HG2	2:B:13:HIS:NE2	2.35	0.42
1:A:218:GLN:HG2	1:A:223:ASP:CA	2.48	0.42
2:B:23:LEU:HD23	2:B:39:LEU:HD22	2.01	0.42
1:A:97:ARG:NH2	1:A:116:TYR:HE1	2.18	0.42
1:A:123:TYR:CZ	1:A:140:ALA:HA	2.55	0.41
1:A:184:ALA:HB1	1:A:266:LEU:HD12	2.02	0.41
2:B:1:ILE:HA	2:B:1:ILE:HD13	1.73	0.41
1:A:166:GLU:OE1	1:A:169:ARG:NH1	2.54	0.41
2:B:36:GLU:H	2:B:36:GLU:HG2	1.74	0.41
1:A:54:GLN:NE2	1:A:174:ASN:HB3	2.36	0.41
1:A:115:GLN:HB2	2:B:60:TRP:HH2	1.86	0.41
1:A:213:ILE:HD12	1:A:263:HIS:HB2	2.03	0.41
2:B:37:VAL:HG13	2:B:82:VAL:HG22	2.03	0.41
2:B:46:ILE:HD13	2:B:46:ILE:HG21	1.74	0.41
1:A:214:THR:HG21	1:A:262:GLN:HE21	1.86	0.40
2:B:9:VAL:HG21	2:B:93:VAL:O	2.21	0.40
1:A:141:GLN:O	1:A:141:GLN:HG3	2.21	0.40
2:B:51:HIS:HA	2:B:65:LEU:O	2.21	0.40
1:A:230:LEU:HD11	1:A:243:LYS:CE	2.52	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	268/270 (99%)	250 (93%)	16 (6%)	2 (1%)	18	38
2	B	97/99 (98%)	95 (98%)	2 (2%)	0	100	100
All	All	365/369 (99%)	345 (94%)	18 (5%)	2 (0%)	24	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	269	PRO
1	A	268	LYS

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	223/226 (99%)	197 (88%)	26 (12%)	5	11
2	B	94/94 (100%)	88 (94%)	6 (6%)	16	35
All	All	317/320 (99%)	285 (90%)	32 (10%)	7	15

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	ILE
1	A	48	ARG
1	A	50	PRO
1	A	55	GLU
1	A	81	LEU
1	A	107	TRP
1	A	115	GLN
1	A	128	GLU
1	A	129	ASP
1	A	141	GLN

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Mol	Chain	Res	Type
1	A	154	GLU
1	A	166	GLU
1	A	173	GLU
1	A	178	THR
1	A	195	SER
1	A	198	GLU
1	A	207	SER
1	A	215	LEU
1	A	226	GLN
1	A	229	GLU
1	A	235	PRO
1	A	249	VAL
1	A	251	SER
1	A	254	GLU
1	A	255	GLN
1	A	266	LEU
2	B	1	ILE
2	B	33	SER
2	B	36	GLU
2	B	40	LEU
2	B	68	THR
2	B	75	LYS

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

Mol	Chain	Res	Type
1	A	114	HIS
1	A	145	HIS
1	A	174	ASN
1	A	262	GLN
2	B	83	ASN

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 [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.