



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:59 PM UTC

PDB ID : 2PAB / pdb_00002pab
Title : STRUCTURE OF PREALBUMIN, SECONDARY, TERTIARY AND QUATERNARY INTERACTIONS DETERMINED BY FOURIER REFINEMENT AT 1.8 ANGSTROMS
Authors : Oatley, S.J.; Blake, C.C.F.
Deposited on : 1977-09-16
Resolution : 1.80 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

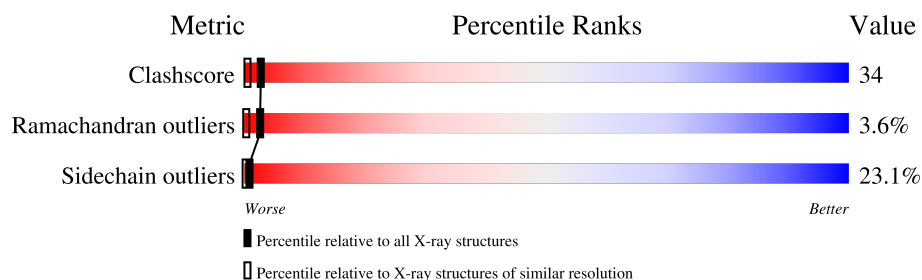
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)

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	127	
1	B	127	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1744 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 TRANSTHYRETIN PRECURSOR.

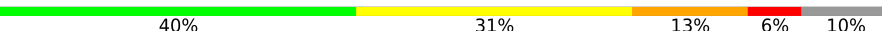
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	114	Total	C	N	O	S	0	0	0
			872	559	142	169	2			
1	B	114	Total	C	N	O	S	0	0	0
			872	559	142	169	2			

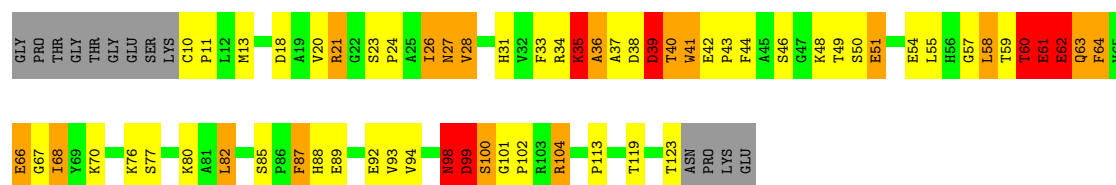
3 Residue-property plots [i](#)

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.

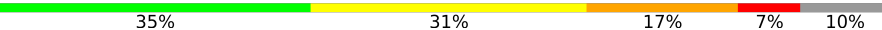
Note EDS was not executed.

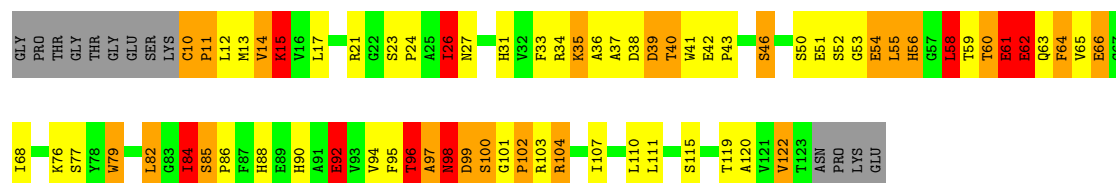
• Molecule 1: TRANSTHYRETIN PRECURSOR

Chain A: 



• Molecule 1: TRANSTHYRETIN PRECURSOR

Chain B: 



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 2	Depositor
Cell constants a, b, c, α , β , γ	43.49Å 85.68Å 65.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 1.80	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	unknown	Depositor
R, R_{free}	0.290 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1744	wwPDB-VP
Average B, all atoms (Å ²)	0.0	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	1.13	3/895 (0.3%)	2.49	55/1221 (4.5%)
1	B	1.12	3/895 (0.3%)	2.62	71/1221 (5.8%)
All	All	1.13	6/1790 (0.3%)	2.55	126/2442 (5.2%)

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	1	0
1	B	2	0
All	All	3	0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	104	ARG	NE-CZ	13.33	1.47	1.33
1	B	34	ARG	NE-CZ	13.14	1.47	1.33
1	A	104	ARG	NE-CZ	13.02	1.47	1.33
1	A	21	ARG	NE-CZ	12.98	1.47	1.33
1	A	34	ARG	NE-CZ	12.90	1.47	1.33
1	B	21	ARG	NE-CZ	12.70	1.47	1.33

All (126) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	63	GLN	N-CA-C	-15.67	94.61	113.88
1	B	99	ASP	CA-CB-CG	-15.38	97.22	112.60
1	B	38	ASP	N-CA-C	15.32	133.08	112.25
1	A	38	ASP	N-CA-C	14.77	129.26	112.57
1	B	40	THR	N-CA-C	14.42	132.49	110.14
1	A	98	ASN	CA-CB-CG	-14.24	98.36	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	40	THR	N-CA-C	13.27	131.00	110.42
1	B	104	ARG	CD-NE-CZ	-12.34	107.12	124.40
1	A	62	GLU	CA-C-N	12.26	144.96	121.54
1	A	62	GLU	C-N-CA	12.26	144.96	121.54
1	A	99	ASP	CA-CB-CG	12.12	124.72	112.60
1	A	36	ALA	N-CA-C	11.15	124.01	110.19
1	B	39	ASP	CA-CB-CG	-10.83	101.77	112.60
1	A	44	PHE	CA-CB-CG	10.72	124.52	113.80
1	A	21	ARG	CD-NE-CZ	-10.39	109.86	124.40
1	B	37	ALA	N-CA-CB	10.31	125.52	110.06
1	B	58	LEU	N-CA-C	9.77	131.61	110.80
1	B	33	PHE	CA-CB-CG	-9.77	104.03	113.80
1	A	62	GLU	O-C-N	9.56	135.30	122.59
1	B	21	ARG	CD-NE-CZ	-9.36	111.30	124.40
1	A	104	ARG	CD-NE-CZ	-9.33	111.33	124.40
1	A	60	THR	N-CA-CB	-9.20	96.69	111.22
1	A	123	THR	N-CA-CB	8.96	126.73	111.50
1	B	40	THR	N-CA-CB	-8.82	97.26	110.85
1	A	102	PRO	N-CA-C	-8.77	94.42	112.47
1	B	27	ASN	CA-CB-CG	-8.72	103.88	112.60
1	A	39	ASP	N-CA-CB	8.59	125.01	110.49
1	B	101	GLY	O-C-N	8.48	130.25	121.77
1	A	58	LEU	N-CA-C	8.46	128.82	110.80
1	B	41	TRP	N-CA-C	8.38	121.91	108.41
1	A	38	ASP	CA-CB-CG	8.38	120.98	112.60
1	B	34	ARG	CD-NE-CZ	-8.33	112.74	124.40
1	A	62	GLU	N-CA-CB	8.31	124.54	110.49
1	B	66	GLU	CB-CA-C	8.29	123.68	109.51
1	B	100	SER	CA-C-N	-8.26	108.89	121.87
1	B	100	SER	C-N-CA	-8.26	108.89	121.87
1	A	61	GLU	CB-CG-CD	8.25	126.62	112.60
1	A	41	TRP	N-CA-C	8.03	121.33	108.41
1	B	41	TRP	CA-C-O	-7.74	112.29	120.58
1	B	15	LYS	CA-CB-CG	-7.69	98.72	114.10
1	A	64	PHE	N-CA-C	-7.61	98.40	109.59
1	A	87	PHE	CA-CB-CG	7.44	121.24	113.80
1	B	63	GLN	CA-C-O	7.42	127.63	119.24
1	B	65	VAL	CA-C-N	-7.29	109.61	122.07
1	B	65	VAL	C-N-CA	-7.29	109.61	122.07
1	B	41	TRP	N-CA-CB	7.16	121.59	110.21
1	B	39	ASP	N-CA-CB	7.13	122.54	110.49
1	A	35	LYS	N-CA-CB	-7.08	98.52	110.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	99	ASP	CA-C-N	7.08	135.07	121.54
1	B	99	ASP	C-N-CA	7.08	135.07	121.54
1	B	39	ASP	CB-CA-C	7.04	124.43	110.42
1	A	39	ASP	CA-C-N	7.01	135.49	122.74
1	A	39	ASP	C-N-CA	7.01	135.49	122.74
1	A	37	ALA	N-CA-CB	7.00	121.00	110.22
1	A	40	THR	CA-CB-CG2	6.92	122.26	110.50
1	A	34	ARG	CD-NE-CZ	-6.84	114.83	124.40
1	B	95	PHE	CA-CB-CG	6.83	120.63	113.80
1	A	61	GLU	N-CA-C	6.78	125.23	110.80
1	A	99	ASP	N-CA-CB	-6.70	99.17	110.49
1	B	34	ARG	N-CA-CB	6.66	121.04	110.65
1	B	64	PHE	N-CA-C	-6.63	99.11	109.25
1	A	37	ALA	N-CA-C	6.61	119.31	111.71
1	B	79	TRP	N-CA-C	6.61	118.14	111.07
1	A	39	ASP	CA-CB-CG	-6.50	106.09	112.60
1	A	39	ASP	CB-CA-C	6.47	123.30	110.42
1	A	101	GLY	O-C-N	6.42	128.19	121.77
1	A	100	SER	N-CA-CB	6.40	121.31	110.49
1	B	38	ASP	CA-CB-CG	6.39	118.99	112.60
1	B	56	HIS	CA-CB-CG	-6.36	107.44	113.80
1	B	92	GLU	CB-CA-C	6.22	121.25	109.37
1	B	63	GLN	CA-CB-CG	-6.21	101.68	114.10
1	A	57	GLY	N-CA-C	6.21	121.29	114.40
1	B	97	ALA	CA-C-N	-6.20	109.70	121.54
1	B	97	ALA	C-N-CA	-6.20	109.70	121.54
1	B	41	TRP	CB-CA-C	-6.20	100.58	110.81
1	B	14	VAL	O-C-N	6.11	129.83	123.05
1	A	40	THR	OG1-CB-CG2	6.10	121.50	109.30
1	A	85	SER	CB-CA-C	-6.07	104.34	110.65
1	B	122	VAL	N-CA-C	6.03	118.05	108.89
1	A	36	ALA	CA-C-N	-6.02	111.61	120.28
1	A	36	ALA	C-N-CA	-6.02	111.61	120.28
1	A	54	GLU	CA-CB-CG	-5.97	102.16	114.10
1	B	42	GLU	N-CA-C	5.93	117.55	110.07
1	B	39	ASP	CA-C-N	5.92	133.34	122.92
1	B	39	ASP	C-N-CA	5.92	133.34	122.92
1	A	35	LYS	CB-CA-C	5.84	122.05	110.42
1	A	61	GLU	CB-CA-C	-5.83	98.82	110.42
1	B	40	THR	CA-C-N	5.83	130.17	122.42
1	B	40	THR	C-N-CA	5.83	130.17	122.42
1	A	104	ARG	N-CA-CB	5.82	120.87	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	103	ARG	CA-C-N	5.81	130.99	121.58
1	B	103	ARG	C-N-CA	5.81	130.99	121.58
1	B	84	ILE	CB-CA-C	-5.74	102.89	110.98
1	B	11	PRO	N-CA-CB	5.72	109.26	103.25
1	B	62	GLU	CB-CA-C	5.69	122.23	112.44
1	B	65	VAL	N-CA-C	-5.67	97.55	109.34
1	B	102	PRO	N-CA-C	-5.59	100.95	112.47
1	B	102	PRO	N-CA-CB	5.58	109.11	103.25
1	A	27	ASN	CA-C-N	-5.53	115.96	122.93
1	A	27	ASN	C-N-CA	-5.53	115.96	122.93
1	A	85	SER	O-C-N	5.52	126.70	120.68
1	B	99	ASP	O-C-N	5.50	129.90	122.59
1	A	66	GLU	CA-CB-CG	-5.46	103.18	114.10
1	B	104	ARG	CA-CB-CG	-5.44	103.22	114.10
1	B	61	GLU	N-CA-C	5.44	122.38	110.80
1	B	26	ILE	CB-CG1-CD1	-5.43	102.39	113.80
1	B	96	THR	CA-C-N	-5.38	115.61	122.77
1	B	96	THR	C-N-CA	-5.38	115.61	122.77
1	B	79	TRP	N-CA-CB	5.35	117.76	110.01
1	B	36	ALA	N-CA-C	5.34	116.81	110.19
1	A	60	THR	CA-C-N	5.27	131.61	121.54
1	A	60	THR	C-N-CA	5.27	131.61	121.54
1	B	88	HIS	CA-CB-CG	-5.23	108.57	113.80
1	B	100	SER	N-CA-C	-5.22	99.68	110.80
1	B	77	SER	O-C-N	5.22	128.10	122.15
1	B	10	CYS	N-CA-C	5.18	125.52	111.00
1	B	76	LYS	O-C-N	5.18	127.48	122.09
1	B	35	LYS	N-CA-CB	-5.16	101.77	110.49
1	A	99	ASP	CB-CA-C	-5.15	100.17	110.42
1	B	111	LEU	N-CA-C	5.13	117.60	109.24
1	B	37	ALA	N-CA-C	5.13	117.54	111.33
1	A	88	HIS	CA-CB-CG	-5.09	108.71	113.80
1	A	63	GLN	CA-CB-CG	-5.07	103.95	114.10
1	A	43	PRO	CA-C-N	-5.05	113.48	122.42
1	A	43	PRO	C-N-CA	-5.05	113.48	122.42
1	B	53	GLY	N-CA-C	-5.02	107.36	115.08

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	58	LEU	CA
1	B	37	ALA	CA

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Mol	Chain	Res	Type	Atom
1	B	39	ASP	CA

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	872	0	841	53	5
1	B	872	0	841	64	8
All	All	1744	0	1682	115	12

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 34.

All (115) 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:98:ASN:ND2	1:A:98:ASN:C	1.79	1.37
1:A:98:ASN:ND2	1:A:99:ASP:N	1.77	1.32
1:A:82:LEU:O	1:A:82:LEU:HD13	1.23	1.29
1:B:97:ALA:O	1:B:98:ASN:CB	1.68	1.29
1:B:61:GLU:O	1:B:62:GLU:OE1	1.53	1.23
1:A:98:ASN:C	1:A:98:ASN:HD22	1.38	1.22
1:B:66:GLU:HB2	1:B:98:ASN:HB3	1.23	1.19
1:A:61:GLU:O	1:A:62:GLU:OE1	1.65	1.12
1:B:79:TRP:CE3	1:B:84:ILE:HD12	1.87	1.10
1:A:82:LEU:O	1:A:82:LEU:CD1	2.08	1.01
1:B:66:GLU:HB2	1:B:98:ASN:CB	1.89	1.00
1:A:26:ILE:HG21	1:A:51:GLU:HA	1.47	0.97
1:A:26:ILE:HG23	1:A:50:SER:O	1.65	0.97
1:B:96:THR:CG2	1:B:98:ASN:HD21	1.78	0.96
1:B:79:TRP:CE3	1:B:84:ILE:CD1	2.50	0.95
1:B:92:GLU:OE2	1:B:94:VAL:HG23	1.66	0.94
1:B:26:ILE:HG22	1:B:50:SER:O	1.68	0.94
1:B:15:LYS:HG3	1:B:54:GLU:HG3	1.49	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:ASN:O	1:B:98:ASN:ND2	2.01	0.93
1:B:79:TRP:HE3	1:B:84:ILE:CD1	1.82	0.92
1:B:97:ALA:O	1:B:98:ASN:CG	2.12	0.91
1:B:96:THR:HG23	1:B:98:ASN:HD21	1.35	0.91
1:A:82:LEU:HD13	1:A:82:LEU:C	1.95	0.91
1:B:96:THR:HG23	1:B:98:ASN:ND2	1.89	0.88
1:B:96:THR:CG2	1:B:98:ASN:ND2	2.37	0.87
1:B:97:ALA:O	1:B:98:ASN:HB3	1.02	0.86
1:B:92:GLU:OE2	1:B:94:VAL:CG2	2.23	0.86
1:B:15:LYS:HD2	1:B:54:GLU:OE1	1.76	0.86
1:A:60:THR:HG22	1:A:63:GLN:HB3	1.62	0.81
1:A:98:ASN:CG	1:A:99:ASP:N	2.39	0.81
1:A:82:LEU:CD1	1:A:82:LEU:C	2.52	0.80
1:A:76:LYS:HE3	1:A:89:GLU:OE2	1.83	0.77
1:B:15:LYS:CG	1:B:54:GLU:HG3	2.15	0.77
1:A:92:GLU:OE2	1:A:94:VAL:CG2	2.34	0.75
1:A:92:GLU:OE2	1:A:94:VAL:HG23	1.87	0.74
1:B:59:THR:OG1	1:B:60:THR:N	2.20	0.74
1:B:66:GLU:HB2	1:B:97:ALA:O	1.88	0.74
1:B:82:LEU:HB3	1:B:84:ILE:HG12	1.69	0.74
1:A:36:ALA:HB2	1:A:42:GLU:HG3	1.69	0.73
1:B:60:THR:HG23	1:B:61:GLU:N	2.02	0.72
1:A:60:THR:HG22	1:A:63:GLN:CB	2.20	0.71
1:A:21:ARG:NH1	1:A:82:LEU:HD21	2.05	0.71
1:A:26:ILE:CG2	1:A:51:GLU:HA	2.21	0.70
1:A:26:ILE:CG2	1:A:50:SER:O	2.40	0.69
1:B:15:LYS:HD2	1:B:54:GLU:HG3	1.74	0.68
1:B:31:HIS:CD2	1:B:46:SER:OG	2.47	0.67
1:B:15:LYS:HD2	1:B:54:GLU:CG	2.24	0.67
1:B:15:LYS:CD	1:B:54:GLU:HG3	2.26	0.65
1:A:98:ASN:HD21	1:A:99:ASP:N	1.91	0.63
1:B:15:LYS:CD	1:B:54:GLU:OE1	2.46	0.63
1:B:79:TRP:CZ3	1:B:84:ILE:HD12	2.33	0.63
1:B:62:GLU:HA	1:B:64:PHE:O	1.98	0.63
1:B:79:TRP:HB3	1:B:84:ILE:HB	1.82	0.62
1:A:98:ASN:ND2	1:A:99:ASP:CA	2.61	0.62
1:A:119:THR:HG23	1:B:115:SER:HB2	1.82	0.62
1:B:60:THR:CG2	1:B:61:GLU:N	2.63	0.61
1:A:35:LYS:HB2	1:A:41:TRP:CZ3	2.37	0.60
1:A:66:GLU:HA	1:A:98:ASN:HB3	1.82	0.60
1:A:18:ASP:OD1	1:A:21:ARG:HD3	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:SER:HB2	1:A:24:PRO:HD2	1.84	0.59
1:B:10:CYS:O	1:B:12:LEU:N	2.37	0.57
1:B:15:LYS:HD2	1:B:54:GLU:CD	2.29	0.57
1:B:92:GLU:OE2	1:B:94:VAL:HG22	2.04	0.57
1:B:13:MET:HG2	1:B:14:VAL:N	2.20	0.56
1:B:12:LEU:HD11	1:B:107:ILE:CD1	2.36	0.56
1:B:82:LEU:HB3	1:B:84:ILE:CG1	2.37	0.55
1:A:76:LYS:CE	1:A:89:GLU:OE2	2.53	0.54
1:A:98:ASN:O	1:A:99:ASP:CB	2.50	0.54
1:B:15:LYS:CE	1:B:54:GLU:OE1	2.56	0.53
1:B:79:TRP:HE3	1:B:84:ILE:CG1	2.22	0.52
1:A:35:LYS:HB3	1:A:68:ILE:O	2.09	0.52
1:A:26:ILE:CG2	1:A:50:SER:C	2.83	0.52
1:B:98:ASN:C	1:B:98:ASN:HD22	2.07	0.51
1:A:92:GLU:HG2	1:A:93:VAL:N	2.26	0.51
1:B:10:CYS:SG	1:B:56:HIS:O	2.61	0.51
1:A:33:PHE:HB2	1:A:70:LYS:HB3	1.93	0.50
1:A:23:SER:HB2	1:A:24:PRO:CD	2.40	0.50
1:B:79:TRP:CZ3	1:B:84:ILE:CD1	2.93	0.50
1:A:28:VAL:HG13	1:A:49:THR:OG1	2.11	0.49
1:A:21:ARG:NH1	1:A:82:LEU:CD2	2.73	0.49
1:A:21:ARG:HH12	1:A:82:LEU:HD21	1.78	0.49
1:B:66:GLU:HB2	1:B:98:ASN:HB2	1.87	0.49
1:B:54:GLU:HB3	1:B:56:HIS:CE1	2.48	0.49
1:B:97:ALA:C	1:B:98:ASN:CG	2.74	0.48
1:A:76:LYS:HB2	1:A:89:GLU:O	2.13	0.48
1:A:92:GLU:OE2	1:A:94:VAL:HG21	2.12	0.47
1:B:55:LEU:HD22	1:B:58:LEU:HD21	1.96	0.47
1:A:99:ASP:HB3	1:A:100:SER:H	1.15	0.47
1:B:12:LEU:HD11	1:B:107:ILE:HD12	1.96	0.47
1:A:36:ALA:HB2	1:A:42:GLU:CG	2.43	0.47
1:A:87:PHE:CE1	1:B:120:ALA:HB2	2.50	0.47
1:B:26:ILE:H	1:B:26:ILE:HG12	1.39	0.47
1:A:66:GLU:HG3	1:A:67:GLY:N	2.29	0.46
1:B:84:ILE:O	1:B:86:PRO:HD3	2.15	0.46
1:A:60:THR:HG23	1:A:61:GLU:N	2.30	0.46
1:B:90:HIS:C	1:B:90:HIS:CD2	2.94	0.45
1:A:62:GLU:HA	1:A:64:PHE:O	2.16	0.45
1:A:18:ASP:OD2	1:A:18:ASP:C	2.60	0.45
1:A:70:LYS:HE3	1:A:92:GLU:HG3	1.98	0.45
1:B:96:THR:HG22	1:B:98:ASN:ND2	2.29	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:ARG:O	1:B:122:VAL:HG13	2.18	0.44
1:A:60:THR:O	1:A:64:PHE:HB2	2.18	0.43
1:A:20:VAL:HG21	1:A:113:PRO:HD3	1.99	0.43
1:B:61:GLU:O	1:B:62:GLU:CD	2.51	0.43
1:A:26:ILE:CG2	1:A:51:GLU:CA	2.94	0.43
1:B:104:ARG:O	1:B:122:VAL:HA	2.18	0.43
1:B:12:LEU:CD1	1:B:107:ILE:HD12	2.50	0.42
1:B:119:THR:HG22	1:B:120:ALA:N	2.35	0.42
1:A:11:PRO:HG2	1:A:59:THR:HG23	2.02	0.42
1:B:82:LEU:HA	1:B:82:LEU:HD22	1.80	0.42
1:B:17:LEU:HB2	1:B:110:LEU:HD12	2.01	0.41
1:A:51:GLU:CA	1:A:51:GLU:OE2	2.68	0.41
1:A:98:ASN:O	1:A:99:ASP:HB2	2.13	0.41
1:B:66:GLU:CB	1:B:98:ASN:CB	2.79	0.41
1:B:23:SER:HB2	1:B:24:PRO:HD2	2.03	0.41

All (12) 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:A:31:HIS:NE2	1:A:62:GLU:CG[4_555]	1.15	1.05
1:B:31:HIS:NE2	1:B:62:GLU:CG[4_456]	1.25	0.95
1:A:31:HIS:NE2	1:A:62:GLU:CB[4_555]	1.44	0.76
1:B:31:HIS:NE2	1:B:62:GLU:CB[4_456]	1.46	0.74
1:B:43:PRO:CG	1:B:102:PRO:CD[4_456]	1.70	0.50
1:B:43:PRO:CB	1:B:102:PRO:CD[4_456]	1.79	0.41
1:B:43:PRO:CG	1:B:102:PRO:CG[4_456]	1.80	0.40
1:A:31:HIS:CD2	1:A:62:GLU:CG[4_555]	1.81	0.39
1:A:31:HIS:CE1	1:A:62:GLU:CB[4_555]	1.88	0.32
1:B:31:HIS:CE1	1:B:62:GLU:CB[4_456]	1.93	0.27
1:B:31:HIS:CD2	1:B:62:GLU:CG[4_456]	1.96	0.24
1:A:82:LEU:O	1:B:85:SER:OG[1_655]	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	112/127 (88%)	101 (90%)	7 (6%)	4 (4%)	2	0
1	B	112/127 (88%)	97 (87%)	11 (10%)	4 (4%)	2	0
All	All	224/254 (88%)	198 (88%)	18 (8%)	8 (4%)	2	0

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	58	LEU
1	A	99	ASP
1	B	98	ASN
1	B	58	LEU
1	B	11	PRO
1	A	39	ASP
1	A	98	ASN
1	B	100	SER

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	93/105 (89%)	71 (76%)	22 (24%)	1	0
1	B	93/105 (89%)	72 (77%)	21 (23%)	1	0
All	All	186/210 (89%)	143 (77%)	43 (23%)	1	0

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	CYS
1	A	13	MET
1	A	26	ILE
1	A	27	ASN
1	A	28	VAL

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Mol	Chain	Res	Type
1	A	35	LYS
1	A	39	ASP
1	A	40	THR
1	A	46	SER
1	A	48	LYS
1	A	51	GLU
1	A	55	LEU
1	A	60	THR
1	A	61	GLU
1	A	62	GLU
1	A	68	ILE
1	A	77	SER
1	A	80	LYS
1	A	82	LEU
1	A	98	ASN
1	A	99	ASP
1	A	104	ARG
1	B	15	LYS
1	B	26	ILE
1	B	35	LYS
1	B	39	ASP
1	B	40	THR
1	B	46	SER
1	B	51	GLU
1	B	52	SER
1	B	54	GLU
1	B	55	LEU
1	B	60	THR
1	B	61	GLU
1	B	62	GLU
1	B	68	ILE
1	B	82	LEU
1	B	84	ILE
1	B	85	SER
1	B	92	GLU
1	B	96	THR
1	B	98	ASN
1	B	99	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	ASN
1	B	31	HIS
1	B	56	HIS
1	B	98	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 [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.