



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 11:27 pm BST

PDB ID : 2PAB
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

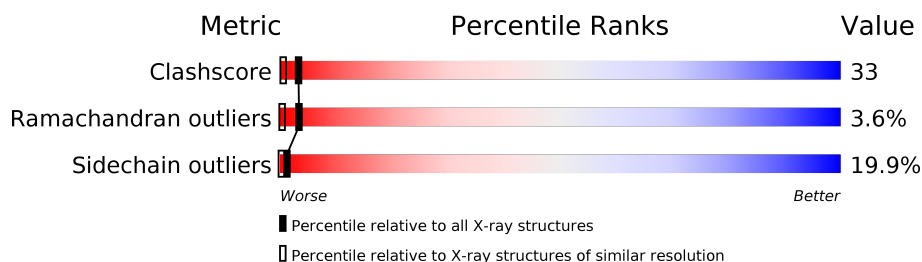
1 Overall quality at a glance

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	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (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 on 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			

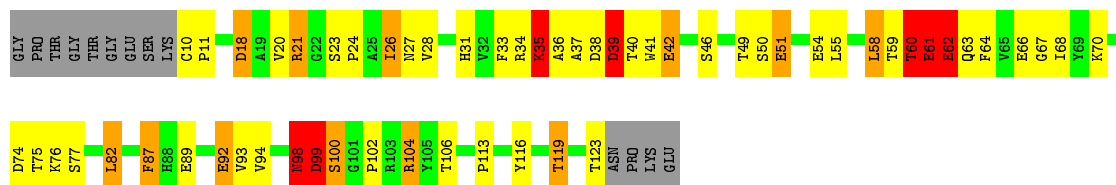
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

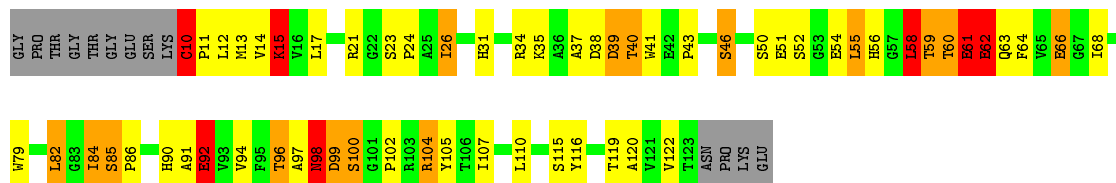
• Molecule 1: TRANSTHYRETIN PRECURSOR

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	0.93	3/895 (0.3%)	2.00	41/1221 (3.4%)
1	B	0.94	3/895 (0.3%)	2.03	33/1221 (2.7%)
All	All	0.93	6/1790 (0.3%)	2.02	74/2442 (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	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	11.22	1.47	1.33
1	B	34	ARG	NE-CZ	11.12	1.47	1.33
1	A	104	ARG	NE-CZ	11.02	1.47	1.33
1	A	21	ARG	NE-CZ	10.98	1.47	1.33
1	A	34	ARG	NE-CZ	10.91	1.47	1.33
1	B	21	ARG	NE-CZ	10.80	1.47	1.33

All (74) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	104	ARG	CD-NE-CZ	-11.79	107.10	123.60
1	B	37	ALA	N-CA-CB	11.02	125.52	110.10
1	A	21	ARG	CD-NE-CZ	-9.82	109.86	123.60
1	A	62	GLU	C-N-CA	9.30	144.96	121.70
1	B	21	ARG	CD-NE-CZ	-8.79	111.29	123.60
1	A	104	ARG	CD-NE-CZ	-8.76	111.33	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	123	THR	N-CA-CB	8.65	126.73	110.30
1	B	38	ASP	N-CA-C	8.17	133.07	111.00
1	A	39	ASP	N-CA-CB	8.01	125.01	110.60
1	B	40	THR	N-CA-C	7.96	132.49	111.00
1	A	62	GLU	O-C-N	7.88	135.31	122.70
1	A	37	ALA	N-CA-CB	7.79	121.00	110.10
1	B	34	ARG	CD-NE-CZ	-7.76	112.74	123.60
1	A	62	GLU	N-CA-CB	7.74	124.54	110.60
1	B	58	LEU	N-CA-C	7.61	131.54	111.00
1	A	40	THR	N-CA-C	7.41	131.00	111.00
1	B	99	ASP	CA-CB-CG	-7.36	97.22	113.40
1	A	62	GLU	CA-C-N	-7.26	101.23	117.20
1	A	100	SER	N-CA-CB	7.21	121.31	110.50
1	A	60	THR	N-CA-CB	-7.17	96.69	110.30
1	A	40	THR	CA-CB-CG2	7.04	122.26	112.40
1	B	39	ASP	CB-CA-C	7.01	124.43	110.40
1	B	40	THR	N-CA-CB	-6.87	97.25	110.30
1	A	98	ASN	CA-CB-CG	-6.84	98.36	113.40
1	A	102	PRO	N-CA-C	-6.80	94.42	112.10
1	A	38	ASP	N-CA-C	6.76	129.26	111.00
1	A	106	THR	CA-CB-CG2	-6.71	103.00	112.40
1	A	35	LYS	N-CA-CB	-6.71	98.52	110.60
1	B	15	LYS	CA-CB-CG	-6.67	98.72	113.40
1	B	66	GLU	CB-CA-C	6.63	123.66	110.40
1	B	39	ASP	N-CA-CB	6.62	122.52	110.60
1	A	58	LEU	N-CA-C	6.59	128.79	111.00
1	B	59	THR	CA-CB-CG2	-6.56	103.22	112.40
1	A	39	ASP	CB-CA-C	6.45	123.30	110.40
1	B	100	SER	C-N-CA	-6.40	108.87	122.30
1	A	99	ASP	N-CA-CB	-6.35	99.17	110.60
1	A	34	ARG	CD-NE-CZ	-6.26	114.83	123.60
1	A	38	ASP	CB-CG-OD2	-6.22	112.70	118.30
1	B	41	TRP	N-CA-CB	6.11	121.59	110.60
1	B	63	GLN	N-CA-C	-6.08	94.57	111.00
1	A	18	ASP	CB-CG-OD1	-6.04	112.86	118.30
1	B	99	ASP	CB-CG-OD2	-5.94	112.95	118.30
1	B	62	GLU	CB-CA-C	5.92	122.23	110.40
1	A	35	LYS	CB-CA-C	5.83	122.05	110.40
1	B	34	ARG	N-CA-CB	5.80	121.04	110.60
1	A	61	GLU	CB-CA-C	-5.79	98.82	110.40
1	A	34	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	A	104	ARG	N-CA-CB	5.71	120.87	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	99	ASP	CA-C-N	-5.58	104.92	117.20
1	A	39	ASP	C-N-CA	5.52	135.49	121.70
1	B	63	GLN	CA-C-N	-5.48	105.15	117.20
1	B	100	SER	N-CA-CB	5.46	118.70	110.50
1	A	75	THR	CA-CB-CG2	-5.43	104.79	112.40
1	B	105	TYR	CB-CG-CD1	-5.42	117.75	121.00
1	B	92	GLU	CB-CA-C	5.42	121.23	110.40
1	A	119	THR	CA-CB-CG2	-5.39	104.85	112.40
1	B	10	CYS	N-CA-C	5.38	125.54	111.00
1	B	99	ASP	C-N-CA	5.35	135.07	121.70
1	A	34	ARG	NE-CZ-NH2	-5.33	117.64	120.30
1	B	63	GLN	CA-CB-CG	-5.33	101.68	113.40
1	B	39	ASP	CA-CB-CG	-5.29	101.77	113.40
1	B	104	ARG	NE-CZ-NH1	-5.28	117.66	120.30
1	A	61	GLU	N-CA-C	5.27	125.23	111.00
1	B	91	ALA	N-CA-CB	5.24	117.43	110.10
1	A	42	GLU	OE1-CD-OE2	5.18	129.51	123.30
1	A	116	TYR	N-CA-CB	5.16	119.88	110.60
1	B	61	GLU	OE1-CD-OE2	5.14	129.47	123.30
1	A	99	ASP	CA-CB-CG	5.14	124.72	113.40
1	B	116	TYR	CB-CG-CD1	-5.13	117.92	121.00
1	A	99	ASP	CB-CA-C	-5.12	100.17	110.40
1	A	54	GLU	CA-CB-CG	-5.11	102.16	113.40
1	A	74	ASP	CB-CG-OD2	-5.08	113.72	118.30
1	A	87	PHE	CB-CG-CD1	-5.07	117.25	120.80
1	A	92	GLU	OE1-CD-OE2	5.03	129.33	123.30

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	58	LEU	CA
1	B	37	ALA	CA
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	840	54	5
1	B	872	0	841	61	8
All	All	1744	0	1681	113	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 33.

All (113) 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:99:ASP:N	1.77	1.32
1:A:82:LEU:HD13	1:A:82:LEU:O	1.23	1.30
1:B:97:ALA:O	1:B:98:ASN:CB	1.68	1.28
1:B:61:GLU:O	1:B:62:GLU:OE1	1.53	1.23
1:A:98:ASN:C	1:A:98:ASN:ND2	1.79	1.21
1:B:66:GLU:HB2	1:B:98:ASN:HB3	1.23	1.18
1:A:98:ASN:C	1:A:98:ASN:HD22	1.38	1.17
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	1.09
1:A:82:LEU:CD1	1:A:82:LEU:O	2.08	1.01
1:B:66:GLU:HB2	1:B:98:ASN:CB	1.89	1.01
1:A:26:ILE:HG21	1:A:51:GLU:HA	1.46	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:15:LYS:HG3	1:B:54:GLU:HG3	1.49	0.94
1:B:26:ILE:HG22	1:B:50:SER:O	1.68	0.94
1:B:92:GLU:OE2	1:B:94:VAL:HG23	1.66	0.94
1:B:98:ASN:ND2	1:B:98:ASN:O	2.01	0.93
1:B:79:TRP:HE3	1:B:84:ILE:CD1	1.82	0.92
1:B:96:THR:HG23	1:B:98:ASN:HD21	1.35	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	0.87
1:B:97:ALA:O	1:B:98:ASN:CG	2.12	0.87
1:A:82:LEU:HD13	1:A:82:LEU:C	1.95	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:B:15:LYS:CG	1:B:54:GLU:HG3	2.15	0.77
1:A:76:LYS:HE3	1:A:89:GLU:OE2	1.83	0.77
1:A:98:ASN:CG	1:A:99:ASP:N	2.39	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:LEU:C	1:A:82:LEU:CD1	2.52	0.75
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:82:LEU:HB3	1:B:84:ILE:HG12	1.69	0.74
1:B:66:GLU:HB2	1:B:97:ALA:O	1.88	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	2.20	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:15:LYS:HD2	1:B:54:GLU:CG	2.24	0.67
1:B:31:HIS:CD2	1:B:46:SER:OG	2.47	0.67
1:B:15:LYS:CD	1:B:54:GLU:HG3	2.26	0.66
1:B:79:TRP:CZ3	1:B:84:ILE:HD12	2.33	0.63
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:62:GLU:HA	1:B:64:PHE:O	1.98	0.62
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	0.62
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
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:92:GLU:OE2	1:B:94:VAL:HG22	2.04	0.57
1:B:13:MET:HG2	1:B:14:VAL:N	2.19	0.56
1:B:12:LEU:HD11	1:B:107:ILE:CD1	2.36	0.56
1:A:98:ASN:O	1:A:99:ASP:CB	2.50	0.55
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:B:15:LYS:CE	1:B:54:GLU:OE1	2.56	0.53
1:B:15:LYS:HD2	1:B:54:GLU:CD	2.29	0.52
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:92:GLU:HG2	1:A:93:VAL:N	2.26	0.51
1:A:33:PHE:HB2	1:A:70:LYS:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:SER:HB2	1:A:24:PRO:CD	2.40	0.50
1:B:10:CYS:SG	1:B:56:HIS:O	2.61	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:54:GLU:HB3	1:B:56:HIS:CE1	2.48	0.49
1:B:66:GLU:HB2	1:B:98:ASN:HB2	1.87	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:79:TRP:CZ3	1:B:84:ILE:CD1	2.93	0.47
1:A:99:ASP:HB3	1:A:100:SER:H	1.15	0.47
1:B:55:LEU:HD22	1:B:58:LEU:HD21	1.96	0.47
1:B:12:LEU:HD11	1:B:107:ILE:HD12	1.96	0.47
1:A:26:ILE:CG2	1:A:50:SER:C	2.82	0.47
1:A:36:ALA:HB2	1:A:42:GLU:CG	2.43	0.47
1:A:66:GLU:HG3	1:A:67:GLY:N	2.29	0.46
1:B:26:ILE:H	1:B:26:ILE:HG12	1.39	0.46
1:A:87:PHE:CE1	1:B:120:ALA:HB2	2.50	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:A:62:GLU:HA	1:A:64:PHE:O	2.16	0.45
1:A:70:LYS:HE3	1:A:92:GLU:HG3	1.98	0.44
1:B:96:THR:HG22	1:B:98:ASN:ND2	2.29	0.44
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:26:ILE:CG2	1:A:51:GLU:CA	2.94	0.43
1:A:20:VAL:HG21	1:A:113:PRO:HD3	1.99	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:B:82:LEU:HA	1:B:82:LEU:HD22	1.80	0.42
1:A:11:PRO:HG2	1:A:59:THR:HG23	2.02	0.42
1:B:17:LEU:HB2	1:B:110:LEU:HD12	2.01	0.42
1:A:51:GLU:CA	1:A:51:GLU:OE2	2.68	0.41
1:B:66:GLU:CB	1:B:98:ASN:CB	2.79	0.41
1:A:98:ASN:O	1:A:99:ASP:HB2	2.13	0.41
1:B:23:SER:HB2	1:B:24:PRO:HD2	2.03	0.41
1:B:90:HIS:C	1:B:90:HIS:CD2	2.94	0.41
1:A:18:ASP:OD2	1:A:18:ASP:C	2.60	0.40
1:A:28:VAL:HG13	1:A:49:THR:HG1	1.86	0.40

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]	1.26	0.94
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]	1.69	0.51
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%)	3	0
1	B	112/127 (88%)	97 (87%)	11 (10%)	4 (4%)	3	0
All	All	224/254 (88%)	198 (88%)	18 (8%)	8 (4%)	3	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

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Mol	Chain	Res	Type
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%)	77 (83%)	16 (17%)	2	0
1	B	93/105 (89%)	72 (77%)	21 (23%)	1	0
All	All	186/210 (89%)	149 (80%)	37 (20%)	1	0

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	CYS
1	A	26	ILE
1	A	27	ASN
1	A	35	LYS
1	A	39	ASP
1	A	46	SER
1	A	51	GLU
1	A	55	LEU
1	A	60	THR
1	A	61	GLU
1	A	62	GLU
1	A	77	SER
1	A	82	LEU
1	A	98	ASN
1	A	99	ASP
1	A	104	ARG
1	B	10	CYS
1	B	15	LYS
1	B	26	ILE
1	B	35	LYS
1	B	39	ASP
1	B	40	THR

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Mol	Chain	Res	Type
1	B	46	SER
1	B	51	GLU
1	B	52	SER
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

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

Mol	Chain	Res	Type
1	A	56	HIS
1	A	98	ASN
1	B	27	ASN
1	B	31	HIS
1	B	56	HIS
1	B	98	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 ⓘ

There are no carbohydrates 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.