



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:13 PM UTC

PDB ID : 1XYL / pdb_00001xyl
Title : THE ROLE OF THE DIVALENT METAL ION IN SUGAR BINDING,
RING OPENING, AND ISOMERIZATION BY D-XYLOSE ISOMERASE:
REPLACEMENT OF A CATALYTIC METAL BY AN AMINO-ACID
Authors : Allen, K.N.; Lavie, A.; Petsko, G.A.; Ringe, D.
Deposited on : 1993-12-07
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

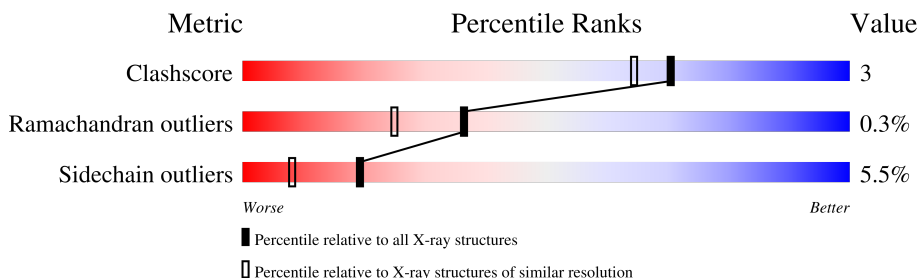
1 Overall quality at a glance

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	386	 80% 18% ..
1	B	386	 80% 16% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8826 atoms, of which 2298 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 XYLOSE ISOMERASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	386	Total	C	H	N	O	S	0	0	0
			3696	1905	672	541	570	8			
1	B	386	Total	C	H	N	O	S	0	0	0
			3696	1905	672	541	570	8			

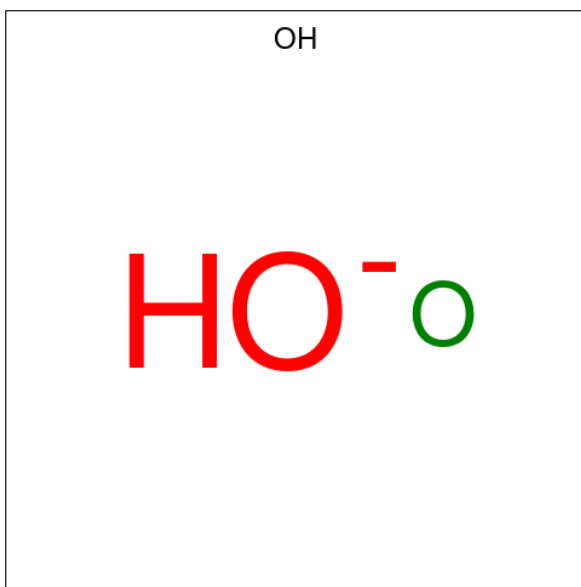
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	180	LYS	GLU	conflict	UNP P15587
B	680	LYS	GLU	conflict	UNP P15587

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

- Molecule 3 is HYDROXIDE ION (CCD ID: OH) (formula: HO).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	H	O	0	0
			2	1	1		
3	B	1	Total	H	O	0	0
			2	1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	238	Total	H	O	0	0
			714	476	238		
4	B	238	Total	H	O	0	0
			714	476	238		

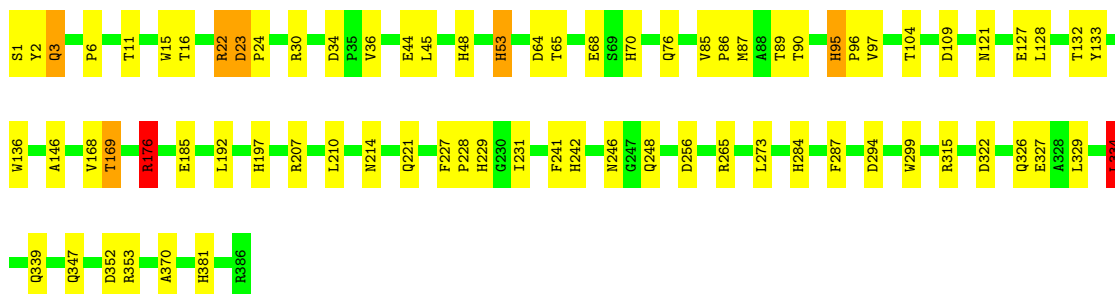
3 Residue-property plots

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.

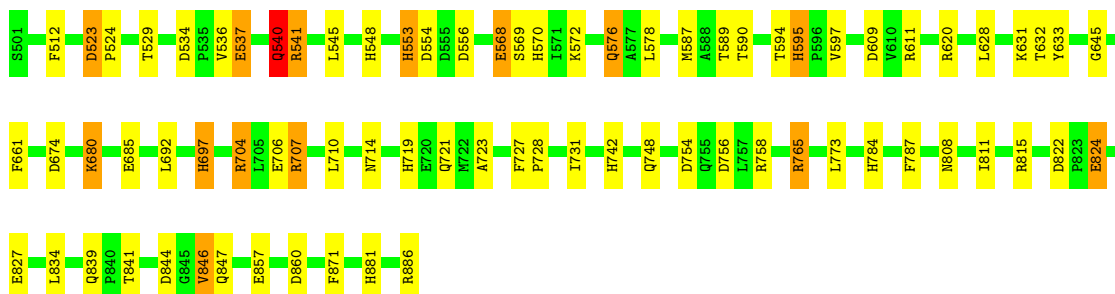
• Molecule 1: XYLOSE ISOMERASE

Chain A: 



• Molecule 1: XYLOSE ISOMERASE

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	87.72Å 99.49Å 94.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.80	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.177 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8826	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, OH

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	14/3096 (0.5%)	1.60	45/4196 (1.1%)
1	B	1.02	11/3096 (0.4%)	1.61	44/4196 (1.0%)
All	All	1.04	25/6192 (0.4%)	1.60	89/8392 (1.1%)

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	B	0	1

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	192	LEU	CA-CB	11.25	1.60	1.52
1	A	48	HIS	CD2-NE2	-7.14	1.29	1.37
1	A	70	HIS	CD2-NE2	-6.59	1.30	1.37
1	A	284	HIS	CD2-NE2	-6.53	1.30	1.37
1	A	95	HIS	CD2-NE2	-6.12	1.31	1.37
1	A	104	THR	CA-CB	6.10	1.61	1.53
1	A	242	HIS	CD2-NE2	-6.06	1.31	1.37
1	B	548	HIS	CD2-NE2	-6.00	1.31	1.37
1	A	381	HIS	CD2-NE2	-5.99	1.31	1.37
1	A	16	THR	CA-CB	5.93	1.61	1.53
1	A	229	HIS	CD2-NE2	-5.90	1.31	1.37
1	B	595	HIS	CD2-NE2	-5.71	1.31	1.37
1	B	595	HIS	CG-ND1	-5.71	1.31	1.38
1	B	570	HIS	CD2-NE2	-5.56	1.31	1.37
1	A	168	VAL	CA-CB	5.50	1.62	1.54
1	A	197	HIS	CD2-NE2	-5.47	1.31	1.37

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	53	HIS	CD2-NE2	-5.40	1.31	1.37
1	B	881	HIS	CD2-NE2	-5.39	1.31	1.37
1	B	742	HIS	CD2-NE2	-5.31	1.32	1.37
1	B	553	HIS	CD2-NE2	-5.27	1.32	1.37
1	B	784	HIS	CG-ND1	-5.19	1.32	1.38
1	B	719	HIS	CD2-NE2	-5.13	1.32	1.37
1	A	284	HIS	CG-ND1	-5.11	1.32	1.38
1	B	692	LEU	CA-CB	5.10	1.56	1.52
1	B	784	HIS	CD2-NE2	-5.04	1.32	1.37

All (89) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	844	ASP	CA-CB-CG	8.76	121.36	112.60
1	B	523	ASP	CA-CB-CG	7.19	119.79	112.60
1	B	609	ASP	CA-CB-CG	7.18	119.78	112.60
1	A	265	ARG	NE-CZ-NH2	-7.15	112.77	119.20
1	A	70	HIS	CA-CB-CG	-7.02	106.78	113.80
1	B	881	HIS	CB-CG-CD2	-6.98	122.13	131.20
1	A	109	ASP	CA-CB-CG	6.88	119.48	112.60
1	B	554	ASP	CA-CB-CG	6.84	119.44	112.60
1	B	576	GLN	OE1-CD-NE2	-6.83	115.77	122.60
1	B	553	HIS	CB-CG-CD2	-6.73	122.45	131.20
1	A	34	ASP	CA-CB-CG	6.62	119.22	112.60
1	B	674	ASP	CA-CB-CG	6.54	119.14	112.60
1	B	645	GLY	N-CA-C	6.43	119.95	112.50
1	A	294	ASP	CA-CB-CG	6.42	119.02	112.60
1	A	242	HIS	CB-CG-CD2	-6.41	122.86	131.20
1	B	824	GLU	CA-CB-CG	-6.29	101.52	114.10
1	A	23	ASP	CA-CB-CG	6.28	118.88	112.60
1	B	754	ASP	CA-CB-CG	6.27	118.87	112.60
1	A	53	HIS	CB-CG-CD2	-6.25	123.08	131.20
1	A	76	GLN	CA-CB-CG	6.22	126.55	114.10
1	B	540	GLN	CB-CA-C	-6.19	100.33	110.85
1	B	765	ARG	NE-CZ-NH1	6.17	127.67	121.50
1	B	822	ASP	CA-C-N	6.16	126.06	119.28
1	B	822	ASP	C-N-CA	6.16	126.06	119.28
1	A	68	GLU	CA-CB-CG	6.16	126.42	114.10
1	A	48	HIS	CB-CG-CD2	-6.16	123.20	131.20
1	B	860	ASP	CA-CB-CG	6.14	118.74	112.60
1	A	381	HIS	CB-CG-CD2	-6.13	123.23	131.20
1	B	537	GLU	N-CA-CB	-6.13	101.11	110.12

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	64	ASP	N-CA-C	6.09	118.72	111.71
1	A	2	TYR	N-CA-C	6.08	123.75	110.80
1	B	808	ASN	OD1-CG-ND2	-5.94	116.66	122.60
1	A	70	HIS	CB-CG-CD2	-5.94	123.47	131.20
1	A	169	THR	N-CA-CB	-5.93	101.37	110.20
1	A	176	ARG	CG-CD-NE	5.91	124.99	112.00
1	B	661	PHE	CA-CB-CG	-5.90	107.90	113.80
1	B	742	HIS	CB-CG-CD2	-5.89	123.54	131.20
1	B	568	GLU	CB-CG-CD	5.88	122.60	112.60
1	B	697	HIS	CB-CG-CD2	-5.73	123.75	131.20
1	A	265	ARG	NE-CZ-NH1	5.71	127.21	121.50
1	B	846	VAL	CG1-CB-CG2	-5.70	98.27	110.80
1	A	322	ASP	CA-CB-CG	5.68	118.28	112.60
1	B	765	ARG	NE-CZ-NH2	-5.67	114.10	119.20
1	A	22	ARG	CB-CG-CD	5.61	124.19	111.30
1	B	556	ASP	CA-CB-CG	5.57	118.17	112.60
1	A	176	ARG	CD-NE-CZ	5.57	132.20	124.40
1	B	723	ALA	N-CA-C	-5.57	106.07	112.92
1	B	871	PHE	CA-CB-CG	5.57	119.37	113.80
1	A	15	TRP	CG-CD2-CE3	5.56	139.46	133.90
1	A	44	GLU	CA-CB-CG	-5.55	102.99	114.10
1	A	246	ASN	OD1-CG-ND2	-5.50	117.10	122.60
1	B	839	GLN	OE1-CD-NE2	-5.50	117.11	122.60
1	A	30	ARG	CA-C-N	5.48	125.38	119.85
1	A	30	ARG	C-N-CA	5.48	125.38	119.85
1	B	534	ASP	CA-CB-CG	5.48	118.08	112.60
1	A	352	ASP	N-CA-C	5.46	117.01	109.15
1	B	537	GLU	CA-CB-CG	5.46	125.01	114.10
1	B	881	HIS	CB-CG-ND1	5.43	130.85	122.70
1	B	594	THR	N-CA-C	5.42	116.99	111.14
1	A	322	ASP	CA-C-N	5.41	125.48	119.32
1	A	322	ASP	C-N-CA	5.41	125.48	119.32
1	A	53	HIS	N-CA-C	-5.38	101.50	110.17
1	A	334	LEU	CA-CB-CG	-5.37	97.51	116.30
1	A	339	GLN	OE1-CD-NE2	-5.34	117.25	122.60
1	A	85	VAL	CA-C-N	5.34	124.95	119.56
1	A	85	VAL	C-N-CA	5.34	124.95	119.56
1	A	214	ASN	CA-C-N	5.33	125.23	119.85
1	A	214	ASN	C-N-CA	5.33	125.23	119.85
1	A	3	GLN	N-CA-C	5.28	116.32	109.72
1	B	569	SER	N-CA-CB	-5.28	102.34	110.16
1	A	207	ARG	CA-C-N	5.13	125.43	119.47

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	207	ARG	C-N-CA	5.13	125.43	119.47
1	A	381	HIS	CA-CB-CG	-5.12	108.68	113.80
1	A	299	TRP	CG-CD2-CE3	5.12	139.02	133.90
1	A	197	HIS	CB-CG-CD2	-5.10	124.57	131.20
1	A	127	GLU	CA-CB-CG	5.08	124.26	114.10
1	B	839	GLN	CA-C-N	5.08	125.08	119.90
1	B	839	GLN	C-N-CA	5.08	125.08	119.90
1	B	512	PHE	CA-CB-CG	5.08	118.88	113.80
1	B	707	ARG	CA-C-N	5.06	124.85	119.28
1	B	707	ARG	C-N-CA	5.06	124.85	119.28
1	B	886	ARG	CA-CB-CG	5.06	124.22	114.10
1	B	534	ASP	CA-C-N	5.05	125.33	119.47
1	B	534	ASP	C-N-CA	5.05	125.33	119.47
1	B	881	HIS	CA-CB-CG	-5.05	108.75	113.80
1	A	136	TRP	CG-CD2-CE3	5.05	138.95	133.90
1	B	822	ASP	CA-CB-CG	5.05	117.65	112.60
1	A	53	HIS	CB-CG-ND1	5.04	130.25	122.70
1	B	590	THR	O-C-N	-5.03	117.31	123.19

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	633	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	3024	672	2916	20	2
1	B	3024	672	2913	22	3
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	1	0	0	0
3	B	1	1	0	0	0
4	A	238	476	0	3	2
4	B	238	476	0	3	3

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6528	2298	5829	40	6

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 3.

All (40) 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:B:537:GLU:HG3	1:B:541:ARG:NH1	2.01	0.76
1:B:595:HIS:HD2	1:B:597:VAL:H	1.37	0.70
1:B:537:GLU:HG3	1:B:541:ARG:HH12	1.56	0.68
1:B:523:ASP:HB2	1:B:524:PRO:HD2	1.75	0.68
1:A:326:GLN:HG3	4:A:1017:HOH:O	1.95	0.67
1:B:765:ARG:HG3	4:B:1248:HOH:O	2.01	0.60
1:A:95:HIS:HD2	1:A:97:VAL:H	1.50	0.60
1:B:680:LYS:HG2	1:B:714:ASN:O	2.02	0.59
1:A:176:ARG:HH11	1:A:176:ARG:HB2	1.69	0.57
1:A:24:PRO:HD3	1:B:524:PRO:HB3	1.89	0.54
1:A:53:HIS:CD2	1:A:89:THR:HG23	2.44	0.53
1:B:595:HIS:CD2	1:B:597:VAL:H	2.23	0.52
1:B:721:GLN:HE21	1:B:748:GLN:HB3	1.76	0.51
1:A:227:PHE:HB3	1:A:228:PRO:HD3	1.92	0.51
1:B:827:GLU:HG3	4:B:1401:HOH:O	2.12	0.50
1:A:176:ARG:HD2	1:A:241:PHE:CE2	2.48	0.49
1:B:704:ARG:HH11	1:B:704:ARG:HB3	1.78	0.49
1:B:756:ASP:HB3	1:B:787:PHE:HA	1.95	0.49
1:B:536:VAL:O	1:B:540:GLN:HG2	2.12	0.48
1:A:227:PHE:CZ	1:A:231:ILE:HD11	2.49	0.47
1:A:347:GLN:HB2	4:A:1458:HOH:O	2.14	0.46
1:B:553:HIS:CD2	1:B:589:THR:HG23	2.51	0.46
1:B:727:PHE:CZ	1:B:731:ILE:HD11	2.51	0.45
1:A:329:LEU:HD23	1:A:334:LEU:HD13	1.97	0.45
1:A:95:HIS:CD2	1:A:97:VAL:H	2.31	0.45
1:A:87:MET:HA	1:A:132:THR:O	2.18	0.43
1:A:221:GLN:HE21	1:A:248:GLN:HB3	1.84	0.43
1:B:727:PHE:HB3	1:B:728:PRO:HD3	2.00	0.43
1:A:256:ASP:HB3	1:A:287:PHE:HA	2.01	0.43
1:A:327:GLU:HG3	4:A:1029:HOH:O	2.18	0.43
1:A:11:THR:HG21	1:A:86:PRO:HG2	2.01	0.42
1:A:121:ASN:HB3	1:A:133:TYR:OH	2.19	0.42
1:B:587:MET:HA	1:B:632:THR:O	2.19	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:ASP:HB2	1:A:24:PRO:HD2	2.00	0.42
1:B:706:GLU:HG3	1:B:707:ARG:HG3	2.02	0.41
1:A:315:ARG:HD3	1:A:315:ARG:HA	1.82	0.41
1:B:811:ILE:O	1:B:815:ARG:HG2	2.21	0.41
1:B:824:GLU:HG2	4:B:1223:HOH:O	2.21	0.40
1:B:540:GLN:HG2	1:B:540:GLN:H	1.78	0.40
1:A:96:PRO:HB3	1:B:529:THR:HG22	2.04	0.40

All (6) 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:B:697:HIS:HE2	4:B:1142:HOH:H1[2_555]	1.00	0.60
1:A:146:ALA:H	4:A:1063:HOH:H1[2_555]	1.15	0.45
1:B:611:ARG:HE	1:B:841:THR:HG1[2_555]	1.15	0.45
4:B:1419:HOH:H1	4:B:1421:HOH:H1[2_555]	1.15	0.45
1:A:370:ALA:H	4:B:1073:HOH:H1[2_555]	1.21	0.39
1:B:758:ARG:H	4:A:1095:HOH:H2[2_555]	1.28	0.32

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	384/386 (100%)	370 (96%)	13 (3%)	1 (0%)	36	25
1	B	384/386 (100%)	371 (97%)	12 (3%)	1 (0%)	36	25
All	All	768/772 (100%)	741 (96%)	25 (3%)	2 (0%)	36	25

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	185	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	685	GLU

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	302/302 (100%)	287 (95%)	15 (5%)	22	10
1	B	302/302 (100%)	284 (94%)	18 (6%)	17	7
All	All	604/604 (100%)	571 (94%)	33 (6%)	19	8

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	SER
1	A	3	GLN
1	A	6	PRO
1	A	22	ARG
1	A	36	VAL
1	A	45	LEU
1	A	65	THR
1	A	90	THR
1	A	128	LEU
1	A	169	THR
1	A	176	ARG
1	A	210	LEU
1	A	273	LEU
1	A	334	LEU
1	A	353	ARG
1	B	540	GLN
1	B	541	ARG
1	B	545	LEU
1	B	568	GLU
1	B	572	LYS
1	B	576	GLN
1	B	578	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	620	ARG
1	B	628	LEU
1	B	631	LYS
1	B	680	LYS
1	B	704	ARG
1	B	710	LEU
1	B	773	LEU
1	B	834	LEU
1	B	846	VAL
1	B	847	GLN
1	B	857	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	53	HIS
1	A	95	HIS
1	A	219	HIS
1	A	221	GLN
1	A	308	ASN
1	A	326	GLN
1	A	376	GLN
1	B	503	GLN
1	B	595	HIS
1	B	621	ASN
1	B	721	GLN
1	B	808	ASN
1	B	876	GLN

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](#)

Of 4 ligands modelled in this entry, 2 are monoatomic and 2 are modelled with single atom - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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.