



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:25 AM UTC

PDB ID : 1PZ2 / pdb_00001pz2
Title : Crystal structure of a transient covalent reaction intermediate of a family 51 alpha-L-arabinofuranosidase
Authors : Hoevel, K.; Shallom, D.; Niefind, K.; Belakhov, V.; Shoham, G.; Baasov, T.; Shoham, Y.; Schomburg, D.
Deposited on : 2003-07-09
Resolution : 2.00 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
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

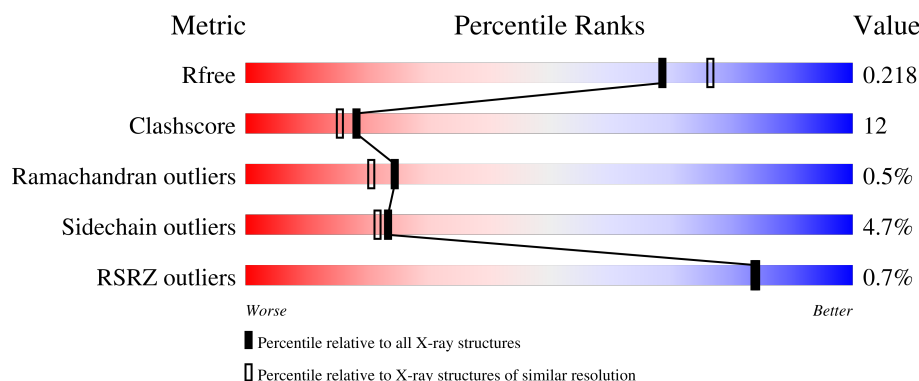
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
R_{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	502	% 80% 16% ..
1	B	502	% 80% 18% ..

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	AHR	A	550	X	-	-	-
2	AHR	B	550	X	-	-	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8757 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.

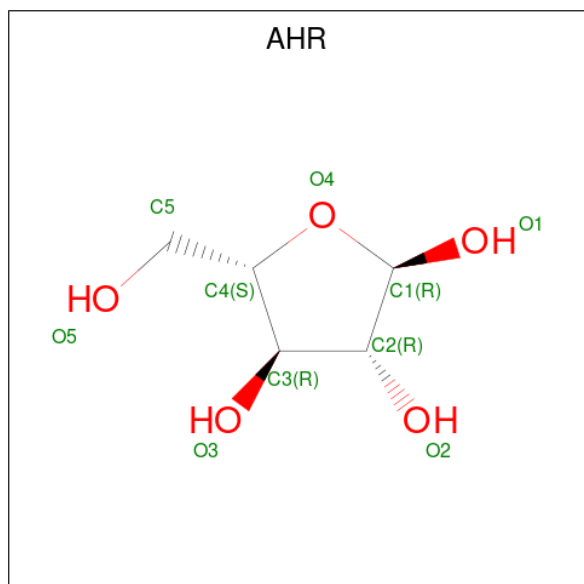
- Molecule 1 is a protein called Alpha-L-arabinofuranosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	497	Total	C	N	O	S	0	0	0
			3986	2540	680	746	20			
1	B	497	Total	C	N	O	S	0	0	0
			3986	2540	680	746	20			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	175	ALA	GLU	engineered mutation	UNP Q9XBQ3
B	175	ALA	GLU	engineered mutation	UNP Q9XBQ3

- Molecule 2 is alpha-L-arabinofuranose (CCD ID: AHR) (formula: C₅H₁₀O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			9	5	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			9	5	4		

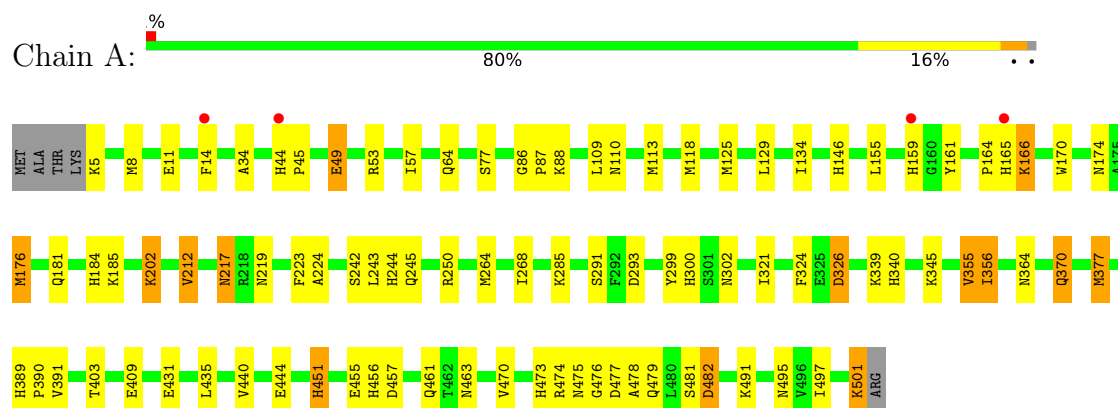
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	399	Total	O	0	0
			399	399		
3	B	368	Total	O	0	0
			368	368		

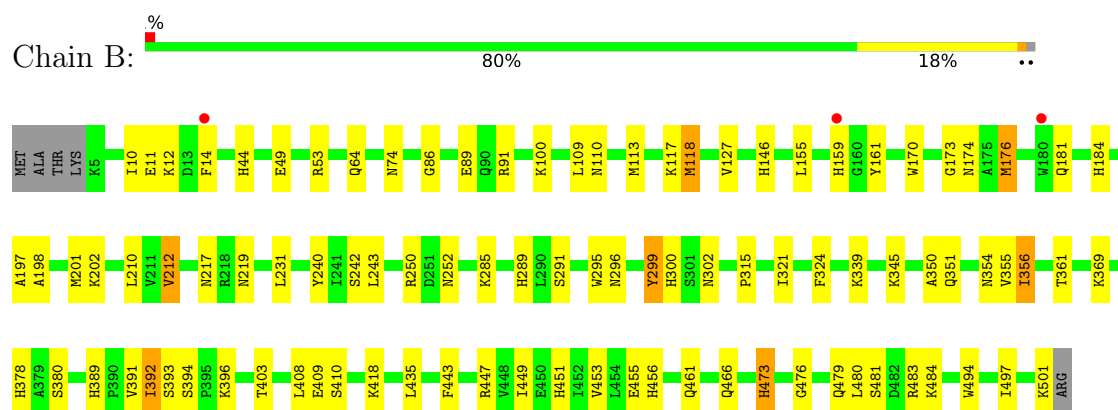
3 Residue-property plots

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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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.

• Molecule 1: Alpha-L-arabinofuranosidase



• Molecule 1: Alpha-L-arabinofuranosidase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	179.43Å 179.43Å 100.23Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.00 20.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	93.3 (20.00-2.00) 93.1 (20.00-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 2.00Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.169 , 0.215 0.179 , 0.218	Depositor DCC
R_{free} test set	3815 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.019 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8757	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	1.06	5/4087 (0.1%)	1.02	2/5553 (0.0%)
1	B	0.97	2/4087 (0.0%)	0.97	5/5553 (0.1%)
All	All	1.02	7/8174 (0.1%)	1.00	7/11106 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	355	VAL	CA-CB	6.89	1.61	1.53
1	A	176	MET	SD-CE	-6.66	1.62	1.79
1	A	377	MET	SD-CE	5.96	1.94	1.79
1	B	176	MET	SD-CE	-5.93	1.64	1.79
1	A	57	ILE	CA-CB	5.70	1.60	1.54
1	A	264	MET	SD-CE	-5.41	1.66	1.79
1	B	91	ARG	C-O	-5.06	1.18	1.24

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	212	VAL	CB-CA-C	-9.05	98.19	111.68
1	B	212	VAL	CB-CA-C	-7.02	101.94	111.63
1	B	380	SER	N-CA-C	6.27	117.78	111.07
1	B	449	ILE	N-CA-C	-5.60	106.25	111.45
1	B	217	ASN	N-CA-CB	-5.32	101.94	110.77
1	B	315	PRO	O-C-N	5.23	123.61	121.15
1	A	326	ASP	N-CA-C	-5.21	105.68	111.36

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	3986	0	3890	106	0
1	B	3986	0	3890	95	0
2	A	9	0	0	0	0
2	B	9	0	0	0	0
3	A	399	0	0	33	1
3	B	368	0	0	25	1
All	All	8757	0	7780	191	2

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 12.

All (191) 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:244:HIS:CE1	3:A:921:HOH:O	1.89	1.20
1:A:44:HIS:ND1	1:A:45:PRO:HD2	1.67	1.07
1:A:391:VAL:HG21	1:B:14:PHE:CZ	1.88	1.07
1:B:451:HIS:ND1	1:B:476:GLY:HA3	1.75	1.01
1:A:340:HIS:CD2	3:A:922:HOH:O	2.13	1.00
1:A:340:HIS:CG	3:A:922:HOH:O	2.19	0.94
1:A:14:PHE:CE2	1:B:391:VAL:HG21	2.04	0.92
1:A:113:MET:HG3	1:A:165:HIS:ND1	1.89	0.87
1:A:451:HIS:CE1	1:A:495:ASN:HD22	1.96	0.84
1:A:155:LEU:O	1:A:159:HIS:HD2	1.61	0.83
1:A:451:HIS:HE1	1:A:495:ASN:HD22	1.26	0.83
1:A:391:VAL:CG2	1:B:14:PHE:CZ	2.62	0.82
1:A:391:VAL:CG2	1:B:14:PHE:CE2	2.64	0.80
1:B:299:TYR:CE1	1:B:300:HIS:CE1	2.68	0.80
1:A:86:GLY:HA2	1:A:159:HIS:ND1	1.97	0.79
1:A:118:MET:HE3	3:A:816:HOH:O	1.83	0.79
1:A:377:MET:SD	3:A:920:HOH:O	2.41	0.79
1:B:410:SER:HB3	3:B:778:HOH:O	1.82	0.78
1:B:456:HIS:CE1	1:B:461:GLN:HG2	2.19	0.78
1:B:86:GLY:N	1:B:159:HIS:HD2	1.83	0.77
1:B:250:ARG:HH21	1:B:302:ASN:HD21	1.33	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:86:GLY:N	1:B:159:HIS:CD2	2.54	0.76
1:A:250:ARG:HH21	1:A:302:ASN:HD21	1.31	0.75
1:A:268:ILE:CG2	1:A:340:HIS:HD2	1.98	0.75
1:A:14:PHE:CZ	1:B:391:VAL:HG21	2.20	0.75
1:A:14:PHE:CE2	1:B:391:VAL:CG2	2.68	0.75
1:B:451:HIS:ND1	1:B:476:GLY:CA	2.49	0.75
1:A:217:ASN:HD21	1:A:219:ASN:HB2	1.51	0.75
1:B:174:ASN:C	3:B:911:HOH:O	2.29	0.74
1:A:86:GLY:CA	1:A:159:HIS:ND1	2.50	0.74
1:B:86:GLY:CA	1:B:159:HIS:CD2	2.71	0.74
1:B:354:ASN:HD21	1:B:361:THR:H	1.32	0.74
1:A:391:VAL:HG21	1:B:14:PHE:CE2	2.23	0.73
1:A:268:ILE:HG21	1:A:340:HIS:HD2	1.54	0.71
1:B:86:GLY:HA2	1:B:159:HIS:CD2	2.25	0.71
1:A:456:HIS:CE1	1:A:461:GLN:HG2	2.28	0.69
1:A:300:HIS:HE1	1:A:326:ASP:OD2	1.76	0.68
1:B:299:TYR:CE1	1:B:300:HIS:HE1	2.10	0.67
1:B:451:HIS:CD2	1:B:497:ILE:HG12	2.28	0.67
1:A:11:GLU:HG2	1:A:14:PHE:CD1	2.30	0.67
1:A:245:GLN:NE2	3:A:789:HOH:O	2.28	0.66
1:B:456:HIS:HE1	1:B:461:GLN:HG2	1.59	0.66
1:B:11:GLU:CG	1:B:14:PHE:HD1	2.08	0.65
1:B:378:HIS:HD2	1:B:494:TRP:HE1	1.45	0.65
1:A:174:ASN:HD22	1:A:181:GLN:HE22	1.45	0.65
1:B:403:THR:HG23	3:B:916:HOH:O	1.97	0.64
1:B:389:HIS:HD2	3:B:773:HOH:O	1.79	0.64
1:A:155:LEU:O	1:A:159:HIS:CD2	2.49	0.64
1:A:268:ILE:HG21	1:A:340:HIS:CD2	2.34	0.63
1:A:11:GLU:HG2	1:A:14:PHE:HD1	1.64	0.62
1:B:299:TYR:CZ	1:B:300:HIS:CE1	2.88	0.61
1:B:339:LYS:NZ	1:B:409:GLU:OE2	2.31	0.61
1:B:11:GLU:HG2	1:B:14:PHE:HD1	1.64	0.61
1:B:86:GLY:H	1:B:159:HIS:HD2	1.49	0.61
1:A:377:MET:HE2	3:A:920:HOH:O	2.00	0.60
1:A:451:HIS:HD2	1:A:478:ALA:HB2	1.65	0.60
1:B:74:ASN:HA	1:B:181:GLN:HE22	1.65	0.60
1:A:155:LEU:HD11	1:A:159:HIS:HE2	1.66	0.60
1:A:44:HIS:ND1	1:A:45:PRO:CD	2.56	0.60
1:B:146:HIS:HD2	3:B:646:HOH:O	1.86	0.59
1:A:44:HIS:HE1	3:A:657:HOH:O	1.86	0.59
1:B:11:GLU:HG2	1:B:14:PHE:CD1	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:451:HIS:HD2	1:B:497:ILE:HG12	1.68	0.58
1:B:219:ASN:ND2	3:B:886:HOH:O	2.36	0.58
1:B:155:LEU:HG	1:B:159:HIS:CE1	2.38	0.58
1:B:174:ASN:HD22	1:B:181:GLN:NE2	2.02	0.58
1:A:244:HIS:ND1	3:A:921:HOH:O	2.12	0.57
1:A:474:ARG:HD3	3:A:841:HOH:O	2.05	0.57
1:B:174:ASN:HD22	1:B:181:GLN:HE22	1.52	0.57
1:A:244:HIS:NE2	1:A:293:ASP:OD1	2.39	0.56
1:A:389:HIS:CD2	3:A:791:HOH:O	2.60	0.55
1:B:378:HIS:HE1	3:B:707:HOH:O	1.88	0.55
1:A:11:GLU:CG	1:A:14:PHE:HD1	2.18	0.55
1:A:377:MET:CE	3:A:920:HOH:O	2.54	0.55
1:A:455:GLU:OE2	1:A:473:HIS:HE1	1.90	0.55
1:A:245:GLN:NE2	3:A:773:HOH:O	2.39	0.55
1:A:268:ILE:CG2	1:A:340:HIS:CD2	2.86	0.55
1:B:202:LYS:HD2	3:B:764:HOH:O	2.06	0.55
1:A:390:PRO:HG3	3:A:887:HOH:O	2.06	0.54
1:A:391:VAL:HG22	1:B:14:PHE:CE2	2.42	0.54
1:A:477:ASP:OD1	1:A:477:ASP:C	2.50	0.54
1:A:49:GLU:HG2	3:A:822:HOH:O	2.06	0.54
1:A:391:VAL:HG21	1:B:14:PHE:CE1	2.41	0.54
1:A:86:GLY:CA	1:A:159:HIS:CE1	2.91	0.54
1:A:390:PRO:HA	3:A:855:HOH:O	2.07	0.54
1:B:455:GLU:OE2	1:B:473:HIS:HE1	1.91	0.53
1:A:185:LYS:HE2	3:A:571:HOH:O	2.09	0.53
1:B:300:HIS:HD2	1:B:321:ILE:O	1.91	0.53
1:B:146:HIS:HE1	3:B:691:HOH:O	1.92	0.52
1:A:377:MET:HG3	3:A:920:HOH:O	2.09	0.52
1:B:453:VAL:HB	1:B:473:HIS:NE2	2.25	0.52
1:A:440:VAL:HG23	1:A:440:VAL:O	2.10	0.52
1:A:340:HIS:CE1	3:A:922:HOH:O	2.61	0.51
1:B:197:ALA:O	1:B:201:MET:HG3	2.10	0.51
1:A:146:HIS:HE1	3:A:687:HOH:O	1.92	0.51
1:A:34:ALA:HB1	3:A:943:HOH:O	2.11	0.51
1:A:501:LYS:HB3	1:A:501:LYS:NZ	2.25	0.51
1:A:364:ASN:CG	3:A:809:HOH:O	2.53	0.51
1:A:321:ILE:HA	1:A:370:GLN:HE22	1.75	0.51
1:B:64:GLN:NE2	3:B:902:HOH:O	2.28	0.51
1:A:86:GLY:HA3	1:A:159:HIS:CE1	2.46	0.50
1:A:451:HIS:HE1	1:A:495:ASN:ND2	2.03	0.50
1:B:392:ILE:HD13	1:B:408:LEU:HD23	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:LEU:O	1:A:113:MET:HG2	2.12	0.50
1:A:155:LEU:CD1	1:A:159:HIS:HE2	2.24	0.50
1:B:453:VAL:HB	1:B:473:HIS:CD2	2.47	0.50
1:A:324:PHE:CZ	1:A:455:GLU:HA	2.46	0.49
1:A:451:HIS:ND1	1:A:476:GLY:HA3	2.27	0.49
1:A:451:HIS:CD2	1:A:497:ILE:HG12	2.48	0.49
1:A:463:ASN:HD21	1:A:470:VAL:H	1.58	0.49
1:B:110:ASN:HB3	1:B:161:TYR:CZ	2.48	0.49
1:B:392:ILE:CD1	1:B:394:SER:HB2	2.42	0.49
1:B:109:LEU:O	1:B:113:MET:HG2	2.13	0.49
1:B:295:TRP:O	1:B:296:ASN:HB2	2.11	0.49
1:A:113:MET:HG3	1:A:165:HIS:CE1	2.47	0.49
1:B:378:HIS:CD2	1:B:494:TRP:HE1	2.27	0.48
1:B:110:ASN:HB3	1:B:161:TYR:CE1	2.48	0.48
1:A:176:MET:HG2	1:A:181:GLN:HG2	1.95	0.48
1:A:473:HIS:HD2	1:A:475:ASN:H	1.61	0.48
1:A:244:HIS:HE1	3:A:643:HOH:O	1.96	0.47
1:B:174:ASN:HB3	3:B:911:HOH:O	2.14	0.47
1:A:364:ASN:HB3	3:A:809:HOH:O	2.14	0.47
1:B:455:GLU:OE2	1:B:473:HIS:CE1	2.68	0.47
1:B:176:MET:O	1:B:184:HIS:CD2	2.68	0.47
1:A:176:MET:O	1:A:184:HIS:HD2	1.98	0.47
1:B:369:LYS:NZ	3:B:867:HOH:O	2.47	0.47
1:B:392:ILE:HG21	3:B:778:HOH:O	2.13	0.47
1:B:117:LYS:HD3	3:B:819:HOH:O	2.15	0.47
1:B:202:LYS:HE3	1:B:210:LEU:HD12	1.97	0.47
1:B:250:ARG:HH21	1:B:302:ASN:ND2	2.07	0.47
1:B:389:HIS:CD2	3:B:773:HOH:O	2.62	0.47
1:B:451:HIS:CG	1:B:476:GLY:HA3	2.49	0.46
1:A:364:ASN:ND2	3:A:842:HOH:O	2.47	0.46
1:B:181:GLN:NE2	3:B:911:HOH:O	2.48	0.46
1:A:118:MET:HA	1:A:118:MET:HE2	1.98	0.46
1:A:110:ASN:HB3	1:A:161:TYR:CZ	2.51	0.46
1:B:127:VAL:O	1:B:173:GLY:HA2	2.16	0.46
1:B:392:ILE:HD12	1:B:393:SER:N	2.31	0.46
1:B:89:GLU:HB3	3:B:887:HOH:O	2.15	0.45
1:B:44:HIS:CD2	1:B:53:ARG:CZ	2.99	0.45
1:A:146:HIS:HD2	3:A:584:HOH:O	1.99	0.45
1:B:473:HIS:CD2	1:B:473:HIS:C	2.94	0.45
1:B:174:ASN:O	3:B:911:HOH:O	2.21	0.45
1:B:202:LYS:HG2	3:B:737:HOH:O	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:PHE:O	1:A:224:ALA:HB3	2.17	0.45
1:A:245:GLN:CD	3:A:773:HOH:O	2.58	0.45
1:B:12:LYS:HB2	1:B:443:PHE:CZ	2.52	0.44
1:A:164:PRO:HG2	1:A:166:LYS:HE2	1.99	0.44
1:B:299:TYR:CE1	1:B:300:HIS:ND1	2.85	0.44
1:B:176:MET:HG2	1:B:181:GLN:HG2	2.00	0.44
1:B:350:ALA:HA	1:B:351:GLN:HA	1.85	0.44
1:A:355:VAL:O	1:A:356:ILE:C	2.60	0.44
1:A:166:LYS:HG2	3:A:810:HOH:O	2.17	0.43
1:A:184:HIS:HE1	3:A:793:HOH:O	2.01	0.43
1:B:100:LYS:NZ	3:B:727:HOH:O	2.47	0.43
1:B:202:LYS:CD	3:B:737:HOH:O	2.67	0.43
1:A:456:HIS:HD2	1:A:457:ASP:H	1.65	0.43
1:B:176:MET:HG3	3:B:911:HOH:O	2.18	0.43
1:A:110:ASN:HB3	1:A:161:TYR:CE1	2.53	0.43
1:A:155:LEU:HG	1:A:159:HIS:CD2	2.53	0.43
1:A:44:HIS:CD2	1:A:53:ARG:CZ	3.01	0.43
1:A:174:ASN:HD22	1:A:181:GLN:NE2	2.14	0.43
1:B:242:SER:HA	1:B:291:SER:O	2.19	0.43
1:B:355:VAL:O	1:B:356:ILE:C	2.62	0.42
1:A:64:GLN:NE2	3:A:854:HOH:O	2.52	0.42
1:A:77:SER:OG	1:A:181:GLN:NE2	2.52	0.42
1:B:118:MET:HE2	1:B:118:MET:HA	2.01	0.42
1:A:339:LYS:NZ	3:A:828:HOH:O	2.52	0.42
1:B:240:TYR:HA	1:B:289:HIS:O	2.20	0.42
1:B:324:PHE:CZ	1:B:455:GLU:HA	2.54	0.42
1:A:339:LYS:NZ	1:A:409:GLU:OE2	2.38	0.42
1:A:176:MET:HE3	1:A:176:MET:HB2	1.79	0.42
1:A:345:LYS:NZ	3:A:879:HOH:O	2.53	0.41
1:A:125:MET:O	1:A:170:TRP:HA	2.20	0.41
1:B:354:ASN:ND2	1:B:361:THR:H	2.09	0.41
1:A:86:GLY:N	1:A:159:HIS:ND1	2.68	0.41
1:B:345:LYS:NZ	3:B:719:HOH:O	2.54	0.41
1:A:217:ASN:ND2	3:A:926:HOH:O	2.53	0.41
1:A:242:SER:HA	1:A:291:SER:O	2.21	0.41
1:B:479:GLN:HG3	3:B:831:HOH:O	2.19	0.41
1:A:129:LEU:O	1:A:185:LYS:HE3	2.21	0.41
1:A:134:ILE:HG21	1:A:134:ILE:HD13	1.81	0.41
1:A:501:LYS:NZ	1:A:501:LYS:CB	2.84	0.41
1:B:170:TRP:CD1	1:B:170:TRP:N	2.89	0.41
1:B:198:ALA:HB1	1:B:210:LEU:HD13	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:PHE:CD2	1:B:14:PHE:CE2	3.08	0.41
1:B:231:LEU:HD23	1:B:231:LEU:HA	1.93	0.41
1:B:300:HIS:CD2	1:B:321:ILE:O	2.72	0.41
1:A:202:LYS:HB2	1:A:202:LYS:HE3	1.94	0.40
1:B:418:LYS:NZ	3:B:787:HOH:O	2.47	0.40
1:A:451:HIS:CG	1:A:476:GLY:HA3	2.56	0.40

All (2) 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)
3:A:667:HOH:O	3:A:924:HOH:O[3_665]	2.13	0.07
3:B:791:HOH:O	3:B:914:HOH:O[3_665]	2.19	0.01

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	495/502 (99%)	476 (96%)	16 (3%)	3 (1%)	21	17
1	B	495/502 (99%)	473 (96%)	20 (4%)	2 (0%)	30	27
All	All	990/1004 (99%)	949 (96%)	36 (4%)	5 (0%)	24	21

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	299	TYR
1	B	299	TYR
1	A	482	ASP
1	B	356	ILE
1	A	356	ILE

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	429/433 (99%)	407 (95%)	22 (5%)	21	19
1	B	429/433 (99%)	411 (96%)	18 (4%)	26	25
All	All	858/866 (99%)	818 (95%)	40 (5%)	23	22

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	8	MET
1	A	49	GLU
1	A	87	PRO
1	A	88	LYS
1	A	166	LYS
1	A	202	LYS
1	A	212	VAL
1	A	217	ASN
1	A	243	LEU
1	A	285	LYS
1	A	370	GLN
1	A	403	THR
1	A	431	GLU
1	A	435	LEU
1	A	444	GLU
1	A	451	HIS
1	A	479	GLN
1	A	481	SER
1	A	482	ASP
1	A	491	LYS
1	A	501	LYS
1	B	10	ILE
1	B	49	GLU
1	B	118	MET
1	B	212	VAL
1	B	243	LEU

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Mol	Chain	Res	Type
1	B	252	ASN
1	B	285	LYS
1	B	392	ILE
1	B	396	LYS
1	B	435	LEU
1	B	447	ARG
1	B	466	GLN
1	B	473	HIS
1	B	480	LEU
1	B	481	SER
1	B	483	ARG
1	B	484	LYS
1	B	501	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	ASN
1	A	64	GLN
1	A	90	GLN
1	A	105	ASN
1	A	146	HIS
1	A	181	GLN
1	A	184	HIS
1	A	217	ASN
1	A	237	HIS
1	A	245	GLN
1	A	252	ASN
1	A	300	HIS
1	A	302	ASN
1	A	340	HIS
1	A	370	GLN
1	A	378	HIS
1	A	389	HIS
1	A	451	HIS
1	A	461	GLN
1	A	463	ASN
1	A	473	HIS
1	A	479	GLN
1	A	495	ASN
1	B	46	GLN
1	B	64	GLN

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Mol	Chain	Res	Type
1	B	181	GLN
1	B	184	HIS
1	B	237	HIS
1	B	252	ASN
1	B	300	HIS
1	B	302	ASN
1	B	354	ASN
1	B	364	ASN
1	B	378	HIS
1	B	473	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	AHR	A	550	1	9,9,10	1.21	1 (11%)	11,12,14	3.99	5 (45%)
2	AHR	B	550	1	9,9,10	0.56	0	11,12,14	3.83	4 (36%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	AHR	A	550	1	2/2/3/4	0/2/15/18	0/1/1/1
2	AHR	B	550	1	2/2/3/4	0/2/15/18	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	550	AHR	C1-C2	3.04	1.57	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	550	AHR	O2-C2-C3	10.95	132.99	111.43
2	A	550	AHR	O2-C2-C3	10.60	132.30	111.43
2	A	550	AHR	O3-C3-C2	4.56	122.78	111.97
2	A	550	AHR	O2-C2-C1	4.03	123.42	110.89
2	B	550	AHR	O2-C2-C1	3.84	122.83	110.89
2	B	550	AHR	O3-C3-C2	3.35	119.91	111.97
2	A	550	AHR	O4-C4-C3	3.25	107.63	104.63
2	A	550	AHR	O4-C1-C2	-2.84	100.87	105.76
2	B	550	AHR	O4-C4-C3	2.76	107.19	104.63

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	550	AHR	C2
2	A	550	AHR	C1
2	B	550	AHR	C2
2	B	550	AHR	C1

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	497/502 (99%)	-0.16	4 (0%) 82 82	27, 38, 58, 73	0
1	B	497/502 (99%)	-0.14	3 (0%) 85 85	30, 41, 58, 69	0
All	All	994/1004 (99%)	-0.15	7 (0%) 84 84	27, 40, 58, 73	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	180	TRP	3.2
1	B	14	PHE	3.2
1	A	14	PHE	3.0
1	A	159	HIS	2.6
1	B	159	HIS	2.5
1	A	165	HIS	2.5
1	A	44	HIS	2.2

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	AHR	A	550	9/10	0.79	0.12	38,42,46,46	0
2	AHR	B	550	9/10	0.87	0.10	40,43,46,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.