



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:53 PM UTC

PDB ID : 8PXU / pdb_00008pxu
Title : Targeting extended blood antigens by Akkermansia muciniphila enzymes unveils a missing link for generating universal donor blood
Authors : Jensen, M.; Abou Hachem, M.; Morth, J.P.
Deposited on : 2023-07-24
Resolution : 1.99 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

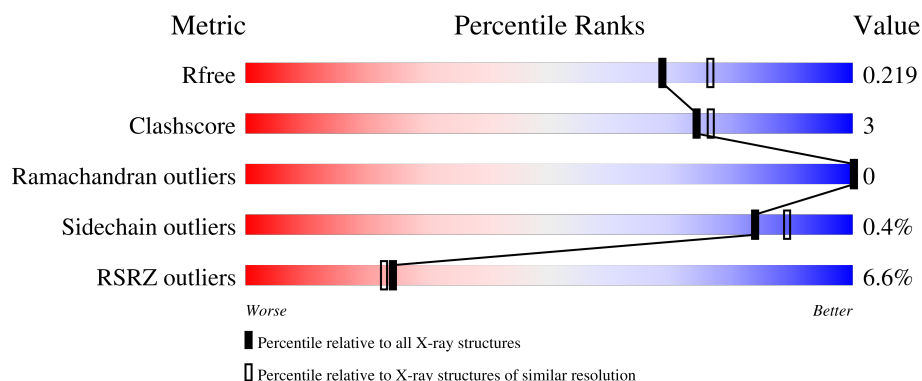
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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	655	6% 91% 7%
1	B	655	9% 89% 8%
1	C	655	4% 89% 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 15945 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 Beta-N-acetylhexosaminidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	646	Total	C	N	O	S	0	0	0
			5079	3228	891	943	17			
1	B	637	Total	C	N	O	S	0	0	0
			5027	3199	880	931	17			
1	C	640	Total	C	N	O	S	8	2	0
			5055	3214	885	938	18			

There are 27 discrepancies between the modelled and reference sequences:

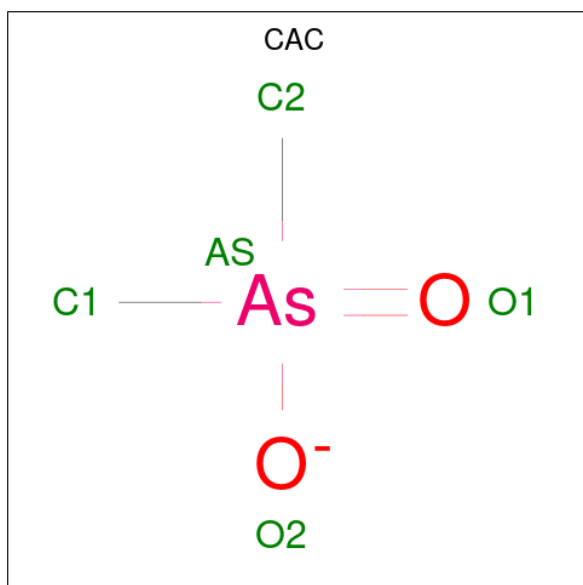
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP B2UN02
A	648	LEU	-	expression tag	UNP B2UN02
A	649	GLU	-	expression tag	UNP B2UN02
A	650	HIS	-	expression tag	UNP B2UN02
A	651	HIS	-	expression tag	UNP B2UN02
A	652	HIS	-	expression tag	UNP B2UN02
A	653	HIS	-	expression tag	UNP B2UN02
A	654	HIS	-	expression tag	UNP B2UN02
A	655	HIS	-	expression tag	UNP B2UN02
B	1	MET	-	initiating methionine	UNP B2UN02
B	648	LEU	-	expression tag	UNP B2UN02
B	649	GLU	-	expression tag	UNP B2UN02
B	650	HIS	-	expression tag	UNP B2UN02
B	651	HIS	-	expression tag	UNP B2UN02
B	652	HIS	-	expression tag	UNP B2UN02
B	653	HIS	-	expression tag	UNP B2UN02
B	654	HIS	-	expression tag	UNP B2UN02
B	655	HIS	-	expression tag	UNP B2UN02
C	1	MET	-	initiating methionine	UNP B2UN02
C	648	LEU	-	expression tag	UNP B2UN02
C	649	GLU	-	expression tag	UNP B2UN02
C	650	HIS	-	expression tag	UNP B2UN02
C	651	HIS	-	expression tag	UNP B2UN02

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Chain	Residue	Modelled	Actual	Comment	Reference
C	652	HIS	-	expression tag	UNP B2UN02
C	653	HIS	-	expression tag	UNP B2UN02
C	654	HIS	-	expression tag	UNP B2UN02
C	655	HIS	-	expression tag	UNP B2UN02

- Molecule 2 is CACODYLATE ION (CCD ID: CAC) (formula: $C_2H_6AsO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	As	C	O	0	0
			5	1	2	2		
2	B	1	Total	As	C	O	0	0
			5	1	2	2		
2	B	1	Total	As	C	O	0	0
			5	1	2	2		
2	C	1	Total	As	C	O	0	0
			5	1	2	2		

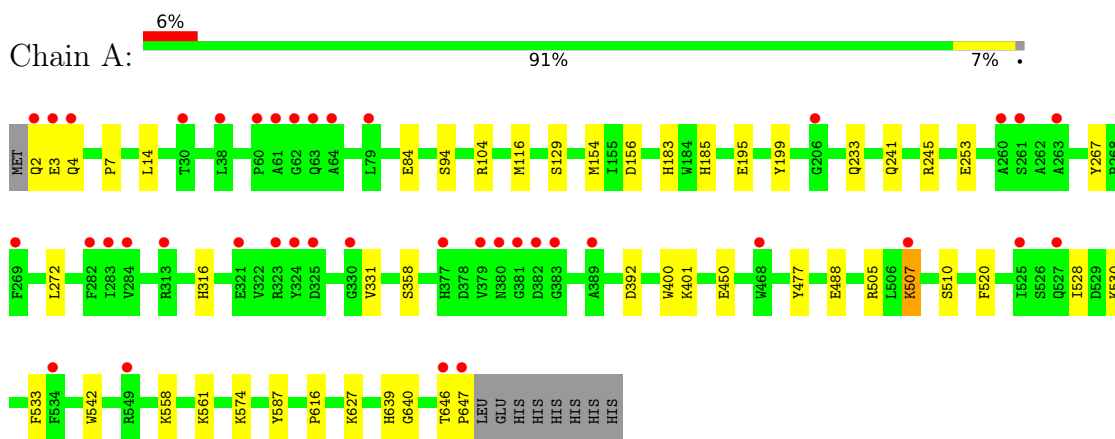
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	253	Total	O	0	0
			253	253		
3	B	237	Total	O	0	0
			237	237		
3	C	274	Total	O	0	0
			274	274		

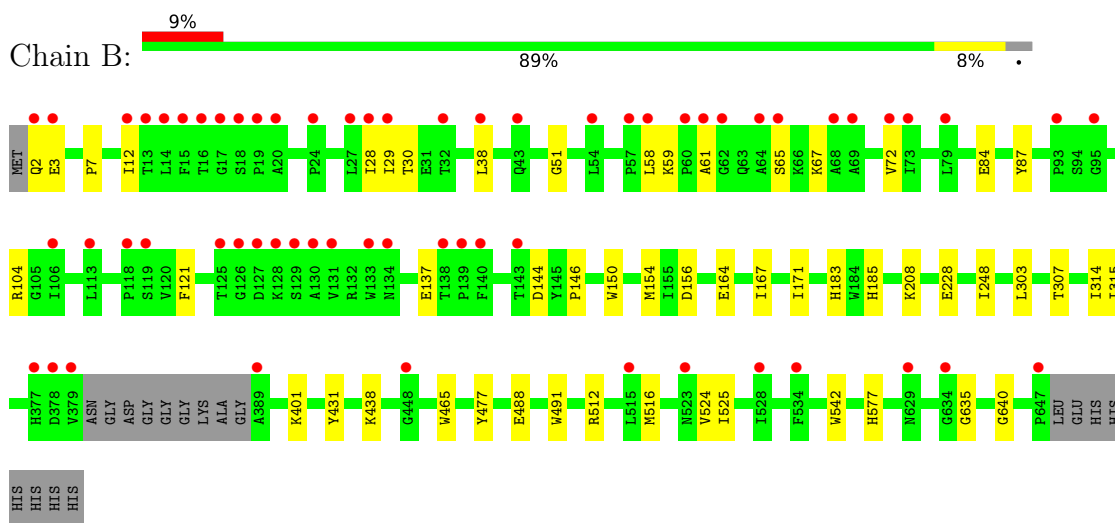
3 Residue-property plots

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.

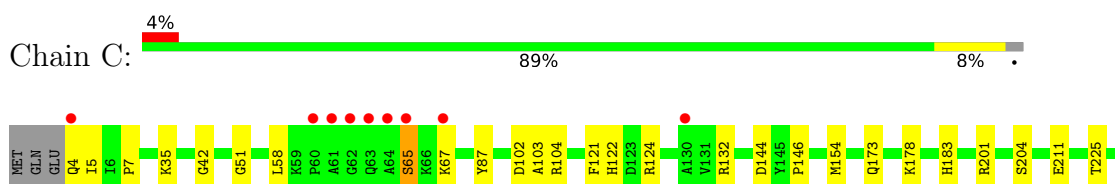
- Molecule 1: Beta-N-acetylhexosaminidase

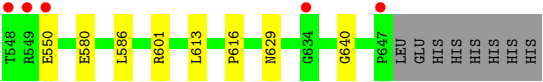
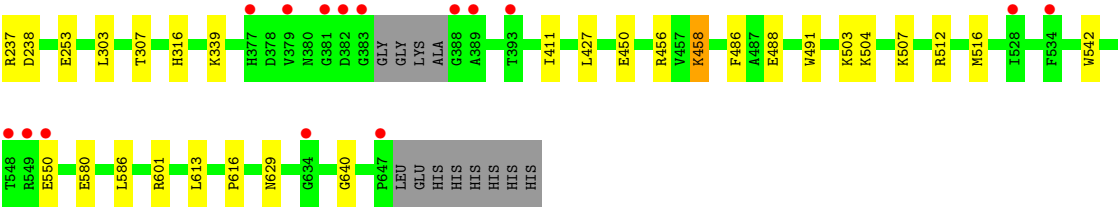


- Molecule 1: Beta-N-acetylhexosaminidase



- Molecule 1: Beta-N-acetylhexosaminidase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	96.71Å 170.74Å 172.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.40 – 1.99 49.40 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.2 (49.40-1.99) 99.3 (49.40-1.99)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.04 (at 1.98Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158, PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.189 , 0.219 0.189 , 0.219	Depositor DCC
R_{free} test set	9731 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	46.3	Xtriage
Anisotropy	0.200	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.009 for -h,l,k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15945	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.29% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.28	0/5207	0.44	0/7049
1	B	0.29	0/5154	0.46	0/6978
1	C	0.29	0/5182	0.45	0/7015
All	All	0.29	0/15543	0.45	0/21042

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	5079	0	5013	28	0
1	B	5027	0	4962	33	0
1	C	5055	0	4985	34	0
2	A	5	0	0	0	0
2	B	10	0	0	0	0
2	C	5	0	0	0	0
3	A	253	0	0	1	0
3	B	237	0	0	2	0
3	C	274	0	0	6	0
All	All	15945	0	14960	92	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 3.

All (92) 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:C:35:LYS:NZ	3:C:802:HOH:O	2.22	0.73
1:B:7:PRO:HB3	1:B:488:GLU:HA	1.77	0.65
1:A:7:PRO:HB3	1:A:488:GLU:HA	1.79	0.65
1:C:104:ARG:NH2	3:C:804:HOH:O	2.35	0.60
1:B:84:GLU:OE1	1:B:104:ARG:NH1	2.38	0.56
1:B:29:ILE:HD12	1:B:61:ALA:HA	1.87	0.56
1:C:542:TRP:CZ2	1:C:640:GLY:HA3	2.40	0.56
1:A:450:GLU:OE2	3:A:901:HOH:O	2.18	0.55
1:C:122:HIS:O	1:C:124:ARG:NH1	2.40	0.54
1:A:84:GLU:OE1	1:A:104:ARG:NH2	2.42	0.53
1:B:156:ASP:HA	1:B:185:HIS:HB3	1.90	0.53
1:C:173:GLN:HE22	1:C:516:MET:HE3	1.74	0.53
1:C:427:LEU:HD21	1:C:486:PHE:HE2	1.74	0.52
1:C:154:MET:HG3	1:C:183:HIS:CG	2.45	0.50
1:C:178:LYS:NZ	3:C:808:HOH:O	2.44	0.50
1:B:307:THR:HG22	1:B:315:ILE:HD11	1.94	0.49
1:C:339:LYS:NZ	3:C:809:HOH:O	2.45	0.49
1:A:272:LEU:HD13	1:A:331:VAL:HG11	1.95	0.49
1:C:173:GLN:HE22	1:C:516:MET:CE	2.25	0.49
1:C:503:LYS:O	1:C:507:LYS:HG3	2.12	0.49
1:C:411:ILE:HG23	1:C:456:ARG:HG3	1.94	0.48
1:A:129:SER:OG	1:C:132:ARG:NH2	2.46	0.48
1:B:2:GLN:HG2	1:B:137:GLU:OE2	2.13	0.48
1:B:2:GLN:HG2	1:B:137:GLU:CD	2.39	0.47
1:A:4:GLN:HB3	1:A:505:ARG:HG2	1.96	0.47
1:C:102:ASP:OD1	1:C:103:ALA:N	2.47	0.47
1:B:65:SER:C	1:B:67:LYS:H	2.22	0.47
1:A:154:MET:HG3	1:A:183:HIS:CG	2.50	0.47
1:A:510:SER:HB2	1:A:520:PHE:CZ	2.50	0.47
1:B:154:MET:HG3	1:B:183:HIS:CG	2.49	0.47
1:C:51:GLY:HA3	1:C:121:PHE:CE2	2.48	0.47
1:A:241:GLN:OE1	1:A:245:ARG:NH1	2.47	0.47
1:B:87:TYR:CE1	1:B:144:ASP:HB3	2.50	0.47
1:B:150:TRP:HZ3	1:B:314:ILE:HD12	1.80	0.47
1:B:150:TRP:CZ3	1:B:314:ILE:HD12	2.50	0.46
1:B:542:TRP:CZ2	1:B:640:GLY:HA3	2.50	0.46
1:A:477:TYR:CE1	1:C:616:PRO:HB3	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:201:ARG:HA	1:C:204:SER:OG	2.16	0.46
1:A:528:ILE:HG23	1:A:533:PHE:HE2	1.81	0.46
1:B:171:ILE:HG23	1:B:248:ILE:HD12	1.97	0.46
1:A:558:LYS:O	1:A:561:LYS:HE2	2.16	0.46
1:A:3:GLU:HG3	1:A:116:MET:HG2	1.97	0.46
1:C:303:LEU:O	1:C:307:THR:HG23	2.16	0.45
1:C:580:GLU:HG3	1:C:601:ARG:HG3	1.98	0.45
1:A:561:LYS:HA	1:A:561:LYS:HD3	1.79	0.45
1:B:3:GLU:HG3	1:B:12:ILE:CD1	2.47	0.45
1:B:438:LYS:NZ	3:B:811:HOH:O	2.49	0.45
1:C:146:PRO:HB3	1:C:491:TRP:CE3	2.52	0.45
1:C:238[B]:ASP:OD1	3:C:801:HOH:O	2.21	0.45
1:C:550:GLU:O	1:C:629:ASN:HA	2.17	0.44
1:A:2:GLN:NE2	1:A:14:LEU:HD13	2.31	0.44
1:A:154:MET:HA	1:A:183:HIS:O	2.17	0.44
1:B:208:LYS:HE3	1:B:228:GLU:OE1	2.17	0.44
1:A:587:TYR:CE1	1:A:627:LYS:HD2	2.52	0.44
1:B:512:ARG:HB3	1:B:516:MET:HE2	2.00	0.44
1:C:87:TYR:CE1	1:C:144:ASP:HB3	2.53	0.44
1:A:156:ASP:HA	1:A:185:HIS:HB3	1.99	0.44
1:A:195:GLU:HG3	1:A:233:GLN:HG3	1.99	0.44
1:B:30:THR:HG21	1:B:38:LEU:HD23	2.00	0.44
1:C:65:SER:O	1:C:65:SER:OG	2.36	0.44
1:B:577:HIS:CE1	1:B:635:GLY:HA3	2.52	0.43
1:B:146:PRO:HB3	1:B:491:TRP:CE3	2.52	0.43
1:A:199:TYR:HB3	1:A:267:TYR:CE2	2.54	0.43
1:A:646:THR:HB	1:A:647:PRO:HD2	2.00	0.43
1:B:28:ILE:HB	1:B:58:LEU:HD23	2.00	0.43
1:B:164:GLU:HA	1:B:167:ILE:HG22	2.00	0.42
1:A:574:LYS:HG2	1:A:639:HIS:HB2	2.01	0.42
1:B:29:ILE:HB	1:B:72:VAL:HG22	2.01	0.42
1:C:237:ARG:NH1	3:C:815:HOH:O	2.51	0.42
1:B:431:TYR:CD1	1:B:524:VAL:HG11	2.53	0.42
1:C:7:PRO:HB3	1:C:488:GLU:HA	2.01	0.42
1:C:458:LYS:HD2	1:C:458:LYS:HA	1.67	0.42
1:C:512:ARG:C	1:C:516:MET:HE2	2.43	0.42
1:C:586:LEU:HB2	1:C:613:LEU:HD13	2.01	0.42
1:A:616:PRO:HB3	1:B:477:TYR:CE1	2.54	0.42
1:B:154:MET:HE1	1:B:465:TRP:CE3	2.54	0.41
1:A:358:SER:OG	1:A:392:ASP:OD2	2.35	0.41
1:A:400:TRP:CE3	1:A:401:LYS:HB2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:GLU:HG3	1:B:12:ILE:HD13	2.03	0.41
1:B:51:GLY:HA3	1:B:121:PHE:CE1	2.55	0.41
1:A:253:GLU:HA	1:A:316:HIS:O	2.21	0.41
1:B:29:ILE:HD13	1:B:59:LYS:O	2.20	0.41
1:C:504:LYS:HD3	1:C:504:LYS:C	2.45	0.41
1:A:542:TRP:CZ2	1:A:640:GLY:HA3	2.56	0.41
1:B:401:LYS:NZ	3:B:819:HOH:O	2.54	0.41
1:C:4:GLN:HB3	1:C:5:ILE:H	1.59	0.41
1:B:303:LEU:O	1:B:307:THR:HG23	2.21	0.41
1:C:42:GLY:HA2	1:C:58:LEU:HD11	2.03	0.41
1:C:211:GLU:HA	1:C:225:THR:O	2.21	0.40
1:A:507:LYS:HB2	1:A:507:LYS:HE3	1.78	0.40
1:B:525:ILE:H	1:B:525:ILE:HD12	1.86	0.40
1:C:253:GLU:HA	1:C:316:HIS:O	2.21	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	644/655 (98%)	631 (98%)	13 (2%)	0	100	100
1	B	633/655 (97%)	612 (97%)	21 (3%)	0	100	100
1	C	638/655 (97%)	616 (97%)	22 (3%)	0	100	100
All	All	1915/1965 (98%)	1859 (97%)	56 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	532/542 (98%)	529 (99%)	3 (1%)	78	85
1	B	528/542 (97%)	528 (100%)	0	100	100
1	C	531/542 (98%)	527 (99%)	4 (1%)	73	80
All	All	1591/1626 (98%)	1584 (100%)	7 (0%)	84	89

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	SER
1	A	507	LYS
1	A	530	LYS
1	C	65	SER
1	C	67	LYS
1	C	450	GLU
1	C	458	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	478	GLN
1	A	610	GLN
1	B	90	GLN
1	B	218	ASN
1	B	241	GLN
1	B	409	ASN
1	B	478	GLN
1	B	577	HIS
1	B	610	GLN
1	C	134	ASN
1	C	173	GLN
1	C	620	GLN
1	C	629	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 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	CAC	A	801	-	2,4,4	3.15	2 (100%)	4,6,6	1.83	1 (25%)
2	CAC	B	702	-	2,4,4	2.47	1 (50%)	4,6,6	1.84	1 (25%)
2	CAC	B	701	-	2,4,4	3.03	2 (100%)	4,6,6	0.96	0
2	CAC	C	701	-	2,4,4	2.57	1 (50%)	4,6,6	1.13	0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	CAC	AS-C1	3.44	1.98	1.90
2	B	702	CAC	AS-C2	3.16	1.97	1.90
2	B	701	CAC	AS-C1	3.12	1.97	1.90
2	C	701	CAC	AS-C1	3.11	1.97	1.90
2	B	701	CAC	AS-C2	2.94	1.97	1.90
2	A	801	CAC	AS-C2	2.82	1.97	1.90

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	801	CAC	O1-AS-C2	-3.43	107.23	111.50
2	B	702	CAC	O2-AS-C2	2.65	112.26	105.84

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

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	646/655 (98%)	0.51	40 (6%)	26 25	37, 55, 81, 117	0
1	B	637/655 (97%)	0.58	62 (9%)	13 12	35, 56, 104, 155	0
1	C	640/655 (97%)	0.31	24 (3%)	44 43	23, 52, 79, 152	2 (0%)
All	All	1923/1965 (97%)	0.47	126 (6%)	24 23	23, 54, 91, 155	2 (0%)

All (126) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	379	VAL	10.1
1	A	647	PRO	7.6
1	C	388	GLY	6.3
1	B	389	ALA	5.8
1	B	64	ALA	5.7
1	C	383	GLY	5.0
1	B	15	PHE	4.7
1	B	61	ALA	4.6
1	C	389	ALA	4.3
1	B	647	PRO	4.2
1	B	131	VAL	4.2
1	B	62	GLY	4.1
1	B	3	GLU	4.1
1	C	549	ARG	4.1
1	C	65	SER	4.0
1	C	61	ALA	4.0
1	B	95	GLY	3.7
1	A	380	ASN	3.7
1	A	64	ALA	3.6
1	C	64	ALA	3.6
1	A	383	GLY	3.6
1	A	284	VAL	3.6
1	A	382	ASP	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	19	PRO	3.5
1	B	65	SER	3.4
1	B	377	HIS	3.4
1	A	206	GLY	3.4
1	B	18	SER	3.4
1	B	14	LEU	3.4
1	B	118	PRO	3.4
1	B	133	TRP	3.3
1	B	129	SER	3.3
1	B	60	PRO	3.3
1	A	379	VAL	3.2
1	A	3	GLU	3.2
1	C	382	ASP	3.2
1	B	138	THR	3.1
1	C	379	VAL	3.1
1	B	126	GLY	3.1
1	B	27	LEU	3.1
1	C	377	HIS	3.1
1	B	634	GLY	3.1
1	B	58	LEU	3.1
1	A	377	HIS	3.1
1	B	106	ILE	3.0
1	B	127	ASP	3.0
1	C	62	GLY	2.9
1	A	527	GLN	2.9
1	B	20	ALA	2.9
1	B	130	ALA	2.9
1	A	325	ASP	2.9
1	B	13	THR	2.8
1	A	61	ALA	2.8
1	C	4	GLN	2.8
1	A	381	GLY	2.8
1	B	128	LYS	2.8
1	C	634	GLY	2.8
1	A	2	GLN	2.7
1	B	73	ILE	2.7
1	B	134	ASN	2.7
1	B	629	ASN	2.7
1	B	125	THR	2.7
1	C	60	PRO	2.7
1	A	525	ILE	2.7
1	B	69	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	381	GLY	2.6
1	B	29	ILE	2.6
1	A	62	GLY	2.6
1	B	378	ASP	2.6
1	C	130	ALA	2.6
1	A	63	GLN	2.6
1	C	528	ILE	2.6
1	A	4	GLN	2.6
1	B	119	SER	2.6
1	C	393	THR	2.6
1	B	16	THR	2.6
1	A	313	ARG	2.5
1	A	534	PHE	2.5
1	B	534	PHE	2.5
1	B	448	GLY	2.5
1	B	57	PRO	2.5
1	C	550	GLU	2.5
1	B	140	PHE	2.5
1	A	646	THR	2.5
1	B	43	GLN	2.5
1	B	17	GLY	2.5
1	A	507	LYS	2.5
1	B	523	ASN	2.4
1	A	260	ALA	2.4
1	A	263	ALA	2.4
1	B	79	LEU	2.4
1	A	549	ARG	2.4
1	B	143	THR	2.4
1	B	139	PRO	2.4
1	B	515	LEU	2.3
1	B	28	ILE	2.3
1	B	54	LEU	2.3
1	C	63	GLN	2.3
1	A	324	TYR	2.3
1	C	548	THR	2.3
1	C	534	PHE	2.3
1	C	647	PRO	2.3
1	A	330	GLY	2.3
1	B	72	VAL	2.3
1	B	2	GLN	2.3
1	B	38	LEU	2.3
1	B	528	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	261	SER	2.3
1	A	38	LEU	2.3
1	C	67	LYS	2.3
1	B	68	ALA	2.2
1	A	389	ALA	2.2
1	A	283	ILE	2.2
1	A	79	LEU	2.2
1	B	24	PRO	2.2
1	B	113	LEU	2.1
1	A	321	GLU	2.1
1	A	282	PHE	2.1
1	A	30	THR	2.1
1	A	468	TRP	2.1
1	A	60	PRO	2.1
1	B	93	PRO	2.1
1	A	323	ARG	2.1
1	B	12	ILE	2.0
1	A	269	PHE	2.0
1	B	32	THR	2.0

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	CAC	A	801	5/5	0.74	0.21	55,62,73,126	0
2	CAC	C	701	5/5	0.83	0.17	60,64,87,123	0
2	CAC	B	702	5/5	0.84	0.18	52,68,86,117	0
2	CAC	B	701	5/5	0.94	0.14	48,62,68,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.