



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:56 PM UTC

PDB ID : 3NA5 / pdb_00003na5
Title : Crystal structure of a bacterial phosphoglucomutase, an enzyme important in the virulence of several human pathogens.
Authors : Mehra-Chaudhary, R.; Beamer, L.J.
Deposited on : 2010-06-01
Resolution : 1.70 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

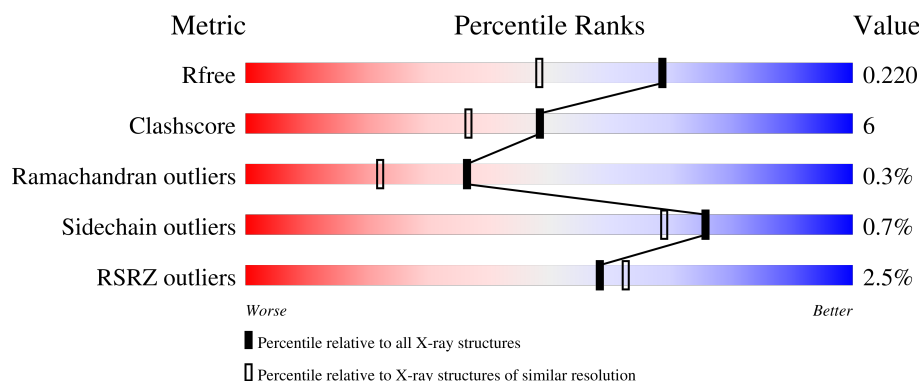
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	570	
1	B	570	

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	BTB	A	600	-	-	X	-
2	BTB	B	600	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9161 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 Phosphoglucomutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	545	Total	C	N	O	S	0	7	0
			4058	2561	706	774	17			
1	B	545	Total	C	N	O	S	0	6	0
			4053	2553	708	775	17			

There are 48 discrepancies between the modelled and reference sequences:

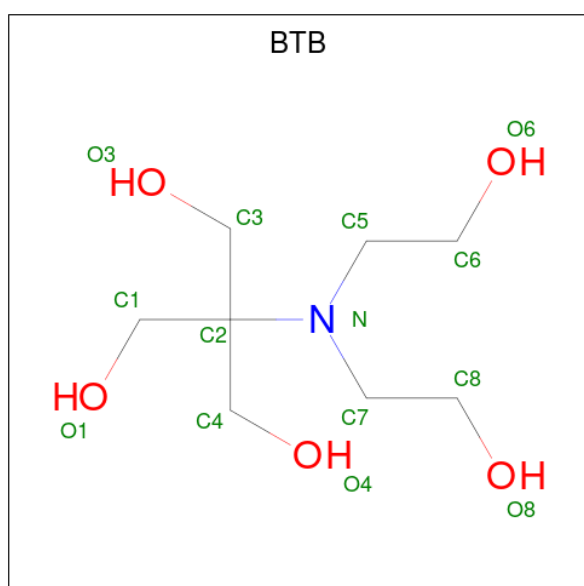
Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	expression tag	UNP Q8ZQW9
A	-22	HIS	-	expression tag	UNP Q8ZQW9
A	-21	HIS	-	expression tag	UNP Q8ZQW9
A	-20	HIS	-	expression tag	UNP Q8ZQW9
A	-19	HIS	-	expression tag	UNP Q8ZQW9
A	-18	HIS	-	expression tag	UNP Q8ZQW9
A	-17	HIS	-	expression tag	UNP Q8ZQW9
A	-16	SER	-	expression tag	UNP Q8ZQW9
A	-15	SER	-	expression tag	UNP Q8ZQW9
A	-14	GLY	-	expression tag	UNP Q8ZQW9
A	-13	VAL	-	expression tag	UNP Q8ZQW9
A	-12	ASP	-	expression tag	UNP Q8ZQW9
A	-11	LEU	-	expression tag	UNP Q8ZQW9
A	-10	GLY	-	expression tag	UNP Q8ZQW9
A	-9	THR	-	expression tag	UNP Q8ZQW9
A	-8	GLU	-	expression tag	UNP Q8ZQW9
A	-7	ASN	-	expression tag	UNP Q8ZQW9
A	-6	LEU	-	expression tag	UNP Q8ZQW9
A	-5	TYR	-	expression tag	UNP Q8ZQW9
A	-4	PHE	-	expression tag	UNP Q8ZQW9
A	-3	GLN	-	expression tag	UNP Q8ZQW9
A	-2	SER	-	expression tag	UNP Q8ZQW9
A	-1	ASN	-	expression tag	UNP Q8ZQW9
A	0	ALA	-	expression tag	UNP Q8ZQW9
B	-23	MET	-	expression tag	UNP Q8ZQW9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	HIS	-	expression tag	UNP Q8ZQW9
B	-21	HIS	-	expression tag	UNP Q8ZQW9
B	-20	HIS	-	expression tag	UNP Q8ZQW9
B	-19	HIS	-	expression tag	UNP Q8ZQW9
B	-18	HIS	-	expression tag	UNP Q8ZQW9
B	-17	HIS	-	expression tag	UNP Q8ZQW9
B	-16	SER	-	expression tag	UNP Q8ZQW9
B	-15	SER	-	expression tag	UNP Q8ZQW9
B	-14	GLY	-	expression tag	UNP Q8ZQW9
B	-13	VAL	-	expression tag	UNP Q8ZQW9
B	-12	ASP	-	expression tag	UNP Q8ZQW9
B	-11	LEU	-	expression tag	UNP Q8ZQW9
B	-10	GLY	-	expression tag	UNP Q8ZQW9
B	-9	THR	-	expression tag	UNP Q8ZQW9
B	-8	GLU	-	expression tag	UNP Q8ZQW9
B	-7	ASN	-	expression tag	UNP Q8ZQW9
B	-6	LEU	-	expression tag	UNP Q8ZQW9
B	-5	TYR	-	expression tag	UNP Q8ZQW9
B	-4	PHE	-	expression tag	UNP Q8ZQW9
B	-3	GLN	-	expression tag	UNP Q8ZQW9
B	-2	SER	-	expression tag	UNP Q8ZQW9
B	-1	ASN	-	expression tag	UNP Q8ZQW9
B	0	ALA	-	expression tag	UNP Q8ZQW9

- Molecule 2 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (CCD ID: BTB) (formula: C₈H₁₉NO₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

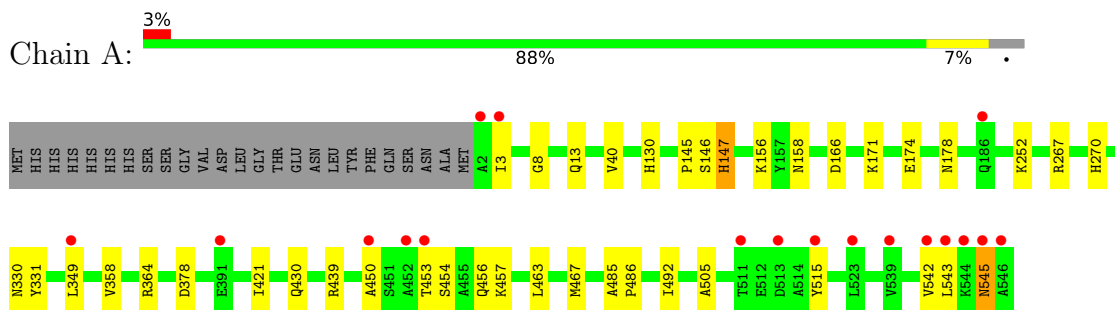
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	502	Total	O	0	0
			502	502		
4	B	518	Total	O	0	0
			518	518		

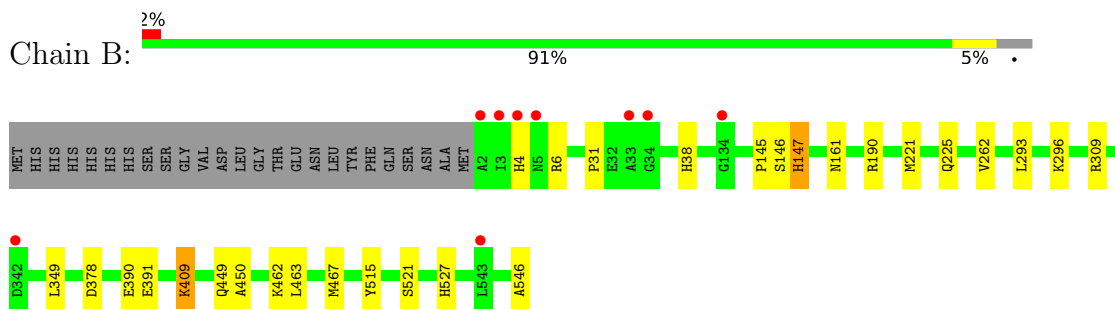
3 Residue-property plots

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: Phosphoglucosyltransferase



• Molecule 1: Phosphoglucosyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.47Å 105.79Å 126.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.27 – 1.70 42.27 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (42.27-1.70) 99.9 (42.27-1.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.72 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.4.0069	Depositor
R, R_{free}	0.182 , 0.218 0.186 , 0.220	Depositor DCC
R_{free} test set	6066 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	27.5	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9161	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.88 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.2117e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: MG, BTB, CSS

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	0.76	0/4158	0.82	0/5657
1	B	0.75	0/4150	0.81	0/5649
All	All	0.75	0/8308	0.81	0/11306

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

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	4058	0	3953	43	0
1	B	4053	0	3936	27	1
2	A	14	0	19	15	0
2	B	14	0	16	13	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	502	0	0	10	1
4	B	518	0	0	7	0
All	All	9161	0	7924	89	1

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 6.

All (89) 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:390:GLU:OE2	1:B:409:LYS:HE2	1.49	1.13
2:A:600:BTB:H62	2:A:600:BTB:H32	1.21	1.07
2:A:600:BTB:C3	2:A:600:BTB:C6	2.38	1.01
2:A:600:BTB:H32	2:A:600:BTB:C6	1.82	1.00
1:A:453:THR:HG23	1:A:456:GLN:OE1	1.60	0.98
2:B:600:BTB:H51	4:B:982:HOH:O	1.63	0.98
1:B:146:SER:OG	1:B:147[A]:HIS:HD2	1.56	0.86
2:A:600:BTB:H41	4:A:984:HOH:O	1.73	0.86
1:A:40:VAL:H	1:A:178:ASN:HD21	1.23	0.85
1:A:349:LEU:HD23	1:A:349:LEU:O	1.82	0.78
2:B:600:BTB:H41	4:B:983:HOH:O	1.86	0.75
1:B:521:SER:OG	1:B:527:HIS:HD2	1.69	0.74
1:B:146:SER:OG	1:B:147[A]:HIS:CD2	2.42	0.72
2:A:600:BTB:C6	2:A:600:BTB:H31	2.18	0.72
1:B:147[A]:HIS:CD2	1:B:147[A]:HIS:H	2.08	0.70
1:A:349:LEU:HD23	1:A:349:LEU:C	2.17	0.69
1:B:378:ASP:OD1	2:B:600:BTB:C4	2.44	0.65
1:A:147[A]:HIS:H	1:A:147[A]:HIS:CD2	2.14	0.65
1:A:146:SER:OG	1:A:147[A]:HIS:HD2	1.80	0.65
1:A:349:LEU:HD21	1:A:505:ALA:CB	2.27	0.64
2:B:600:BTB:C1	2:B:600:BTB:O6	2.44	0.62
1:B:378:ASP:OD1	2:B:600:BTB:H41	2.01	0.60
1:A:147[B]:HIS:CD2	4:A:1033:HOH:O	2.55	0.59
1:B:4:HIS:CD2	1:B:6:ARG:H	2.22	0.58
1:B:463:LEU:HD22	1:B:467:MET:HE2	1.86	0.58
1:A:463:LEU:HD22	1:A:542:VAL:CG1	2.34	0.58
1:B:521:SER:OG	1:B:527:HIS:CD2	2.54	0.58
1:A:453:THR:O	1:A:457:LYS:HG3	2.04	0.57
1:A:349:LEU:HD21	1:A:505:ALA:HB2	1.86	0.57
1:A:40:VAL:N	1:A:178:ASN:HD21	1.97	0.56
1:A:156:LYS:NZ	1:A:158:ASN:HD21	2.03	0.56
2:B:600:BTB:O6	2:B:600:BTB:H11	2.05	0.55
2:A:600:BTB:H62	2:A:600:BTB:C3	2.07	0.55
1:B:463:LEU:HD22	1:B:467:MET:CE	2.36	0.54
1:B:463:LEU:CD2	1:B:467:MET:HE1	2.39	0.53
1:A:378:ASP:OD1	2:A:600:BTB:C4	2.57	0.53
1:B:293:LEU:O	1:B:296:LYS:HG2	2.09	0.53
1:A:166:ASP:HB2	2:A:600:BTB:O1	2.09	0.53
1:B:161:ASN:ND2	2:B:600:BTB:H72	2.24	0.53
1:B:161:ASN:ND2	2:B:600:BTB:H42	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:600:BTB:H61	4:B:847:HOH:O	2.11	0.51
1:A:349:LEU:HD11	1:A:492:ILE:HD12	1.93	0.50
1:B:190:ARG:NH2	4:B:910:HOH:O	2.38	0.50
1:B:221:MET:O	1:B:225:GLN:HG3	2.12	0.50
1:A:364:ARG:NH2	4:A:948:HOH:O	2.39	0.49
1:A:450:ALA:HB3	1:A:515:TYR:CE2	2.48	0.48
1:B:463:LEU:CD2	1:B:467:MET:CE	2.91	0.48
1:A:130:HIS:HE1	4:A:814:HOH:O	1.98	0.47
1:A:463:LEU:CD2	1:A:542:VAL:CG1	2.93	0.47
1:A:13[A]:GLN:NE2	1:B:262:VAL:H	2.13	0.47
2:A:600:BTB:H12	2:A:600:BTB:H71	1.30	0.47
1:A:349:LEU:C	1:A:349:LEU:CD2	2.86	0.46
1:A:378:ASP:OD1	2:A:600:BTB:H41	2.15	0.46
1:A:463:LEU:HD22	1:A:542:VAL:HG11	1.97	0.46
2:B:600:BTB:C3	4:B:983:HOH:O	2.63	0.46
2:B:600:BTB:C6	4:B:847:HOH:O	2.63	0.46
1:A:40:VAL:H	1:A:178:ASN:ND2	2.02	0.46
2:A:600:BTB:N	4:A:984:HOH:O	2.36	0.45
1:B:546:ALA:C	4:B:915:HOH:O	2.59	0.45
1:A:13[A]:GLN:HE22	1:A:267:ARG:HH12	1.65	0.45
1:A:171:LYS:NZ	1:A:174:GLU:OE2	2.48	0.45
2:A:600:BTB:H31	2:A:600:BTB:O6	2.17	0.44
1:B:147[A]:HIS:CE1	1:B:309:ARG:NH1	2.86	0.44
1:A:454:SER:HB3	4:A:988:HOH:O	2.17	0.44
1:A:156:LYS:HZ1	1:A:158:ASN:HD21	1.64	0.44
2:B:600:BTB:H51	2:B:600:BTB:H11	1.50	0.43
1:A:331:TYR:CG	1:A:421:ILE:HG12	2.53	0.43
1:A:146:SER:OG	1:A:147[A]:HIS:CD2	2.67	0.43
1:B:463:LEU:HD23	1:B:467:MET:HE1	2.00	0.43
1:A:166:ASP:OD2	2:A:600:BTB:O3	2.36	0.42
1:A:252:LYS:HE3	4:A:879:HOH:O	2.18	0.42
1:A:330:ASN:HD22	1:A:358:VAL:HG13	1.84	0.42
2:A:600:BTB:H32	4:A:984:HOH:O	2.20	0.42
1:A:463:LEU:HD21	1:A:542:VAL:HG12	2.02	0.42
1:A:430:GLN:NE2	4:A:870:HOH:O	2.53	0.42
1:B:31:PRO:HB3	1:B:38:HIS:HB3	2.01	0.42
2:B:600:BTB:H72	2:B:600:BTB:H42	1.20	0.41
2:A:600:BTB:H51	2:A:600:BTB:H11	1.31	0.41
1:A:467:MET:HE2	1:A:542:VAL:HG13	2.01	0.41
1:B:409:LYS:H	1:B:409:LYS:HG3	1.73	0.41
1:B:449:GLN:HG3	1:B:515:TYR:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:485:ALA:HA	1:A:486:PRO:HD3	1.97	0.41
1:A:8:GLY:O	1:A:270:HIS:HD2	2.04	0.40
1:A:439:ARG:NH2	4:A:803:HOH:O	2.54	0.40
1:B:449:GLN:HG2	1:B:450:ALA:N	2.35	0.40
1:A:543:LEU:C	1:A:545:ASN:N	2.79	0.40

All (1) 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:462:LYS:NZ	4:A:891:HOH:O[4_456]	2.14	0.06

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	549/570 (96%)	540 (98%)	7 (1%)	2 (0%)	30	16
1	B	548/570 (96%)	539 (98%)	8 (2%)	1 (0%)	43	28
All	All	1097/1140 (96%)	1079 (98%)	15 (1%)	3 (0%)	36	22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	545	ASN
1	B	145	PRO
1	A	145	PRO

5.3.2 Protein sidechains [i](#)

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	412/443 (93%)	409 (99%)	3 (1%)	76	69
1	B	412/443 (93%)	407 (99%)	5 (1%)	63	51
All	All	824/886 (93%)	816 (99%)	8 (1%)	76	58

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ILE
1	A	147[A]	HIS
1	A	147[B]	HIS
1	B	147[A]	HIS
1	B	147[B]	HIS
1	B	349	LEU
1	B	391	GLU
1	B	409	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	25	GLN
1	A	54	HIS
1	A	114	ASN
1	A	130	HIS
1	A	158	ASN
1	A	178	ASN
1	A	270	HIS
1	A	330	ASN
1	A	430	GLN
1	B	4	HIS
1	B	25	GLN
1	B	250	HIS
1	B	270	HIS
1	B	430	GLN
1	B	434	ASN
1	B	501	ASN
1	B	527	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	CSS	B	285	1	4,6,7	0.82	0	2,6,8	0.98	0
1	CSS	A	285	1	4,6,7	0.67	0	2,6,8	0.62	0

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
1	CSS	B	285	1	-	0/1/5/7	-
1	CSS	A	285	1	-	0/1/5/7	-

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.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	BTB	A	600	-	13,13,13	0.85	1 (7%)	7,16,16	1.71	2 (28%)
2	BTB	B	600	-	13,13,13	1.16	2 (15%)	7,16,16	2.75	2 (28%)

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	BTB	A	600	-	-	9/21/21/21	-
2	BTB	B	600	-	-	7/21/21/21	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	600	BTB	C3-C2	-2.34	1.50	1.53
2	B	600	BTB	C1-C2	-2.22	1.50	1.53
2	A	600	BTB	C2-N	-2.00	1.44	1.48

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	600	BTB	O1-C1-C2	-5.36	98.79	111.40
2	B	600	BTB	O4-C4-C2	-4.57	100.64	111.40
2	A	600	BTB	O4-C4-C2	-3.29	103.67	111.40
2	A	600	BTB	O1-C1-C2	-2.34	105.91	111.40

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	600	BTB	O1-C1-C2-C4
2	A	600	BTB	O1-C1-C2-N
2	A	600	BTB	C1-C2-C3-O3
2	A	600	BTB	N-C2-C3-O3
2	A	600	BTB	C1-C2-C4-O4
2	A	600	BTB	C3-C2-C4-O4
2	B	600	BTB	O1-C1-C2-C4
2	B	600	BTB	C1-C2-C3-O3
2	B	600	BTB	N-C2-C3-O3
2	B	600	BTB	C1-C2-C4-O4
2	B	600	BTB	C3-C2-C4-O4
2	B	600	BTB	C6-C5-N-C2
2	A	600	BTB	N-C5-C6-O6
2	A	600	BTB	N-C7-C8-O8
2	B	600	BTB	O1-C1-C2-C3
2	A	600	BTB	O1-C1-C2-C3

There are no ring outliers.

2 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	BTB	15	0
2	B	600	BTB	13	0

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	544/570 (95%)	0.04	18 (3%)	49	53	11, 24, 44, 87	9 (1%)
1	B	544/570 (95%)	0.06	9 (1%)	69	73	14, 26, 39, 61	6 (1%)
All	All	1088/1140 (95%)	0.05	27 (2%)	58	62	11, 25, 41, 87	15 (1%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	391	GLU	5.9
1	B	3	ILE	5.6
1	B	2	ALA	4.3
1	A	546	ALA	3.9
1	A	545	ASN	3.7
1	A	542	VAL	3.3
1	A	186	GLN	3.3
1	A	3	ILE	3.1
1	A	349	LEU	3.0
1	A	453	THR	2.8
1	B	33	ALA	2.7
1	A	513	ASP	2.7
1	B	34	GLY	2.7
1	B	134	GLY	2.5
1	B	4	HIS	2.5
1	A	450	ALA	2.5
1	A	452	ALA	2.4
1	B	5	ASN	2.4
1	A	2	ALA	2.4
1	A	515	TYR	2.4
1	A	544	LYS	2.3
1	A	523	LEU	2.2
1	A	511	THR	2.2
1	A	539	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	342	ASP	2.0
1	A	543	LEU	2.0
1	B	543	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	CSS	A	285	7/8	0.83	0.12	24,25,30,32	0
1	CSS	B	285	7/8	0.85	0.12	26,27,33,33	0

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(Å ²)	Q<0.9
2	BTB	A	600	14/14	0.78	0.20	25,32,37,38	14
2	BTB	B	600	14/14	0.84	0.15	21,27,35,35	14
3	MG	A	547	1/1	0.99	0.02	18,18,18,18	0
3	MG	B	547	1/1	0.99	0.03	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.