



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 03:36 AM UTC

PDB ID : 3AAG / pdb_00003aag
Title : Crystal structure of C. jejuni pglb C-terminal domain
Authors : Maita, N.; Kohda, D.
Deposited on : 2009-11-16
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

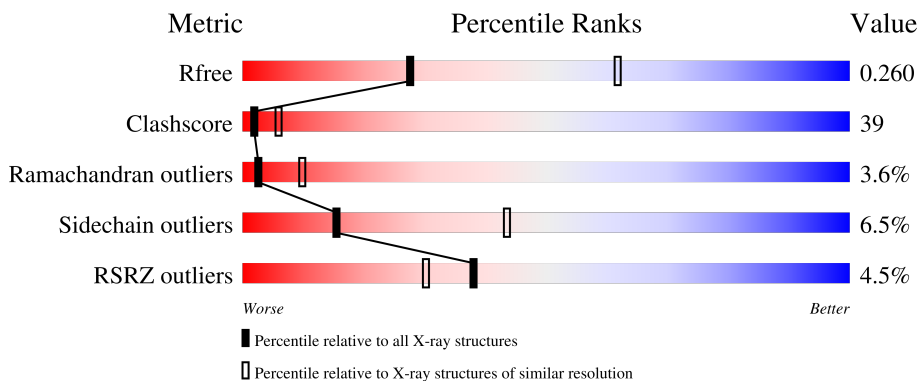
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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 $\leq 5\%$. 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	291	
1	B	291	

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
1	MLY	A	707	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4036 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 General glycosylation pathway protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	267	2176	1426	336	406	8	0	0	0
1	B	236	1859	1221	286	346	6	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	423	GLY	-	expression tag	UNP Q5HTX9
A	424	PRO	-	expression tag	UNP Q5HTX9
A	425	LEU	-	expression tag	UNP Q5HTX9
A	426	GLY	-	expression tag	UNP Q5HTX9
A	427	SER	-	expression tag	UNP Q5HTX9
B	423	GLY	-	expression tag	UNP Q5HTX9
B	424	PRO	-	expression tag	UNP Q5HTX9
B	425	LEU	-	expression tag	UNP Q5HTX9
B	426	GLY	-	expression tag	UNP Q5HTX9
B	427	SER	-	expression tag	UNP Q5HTX9

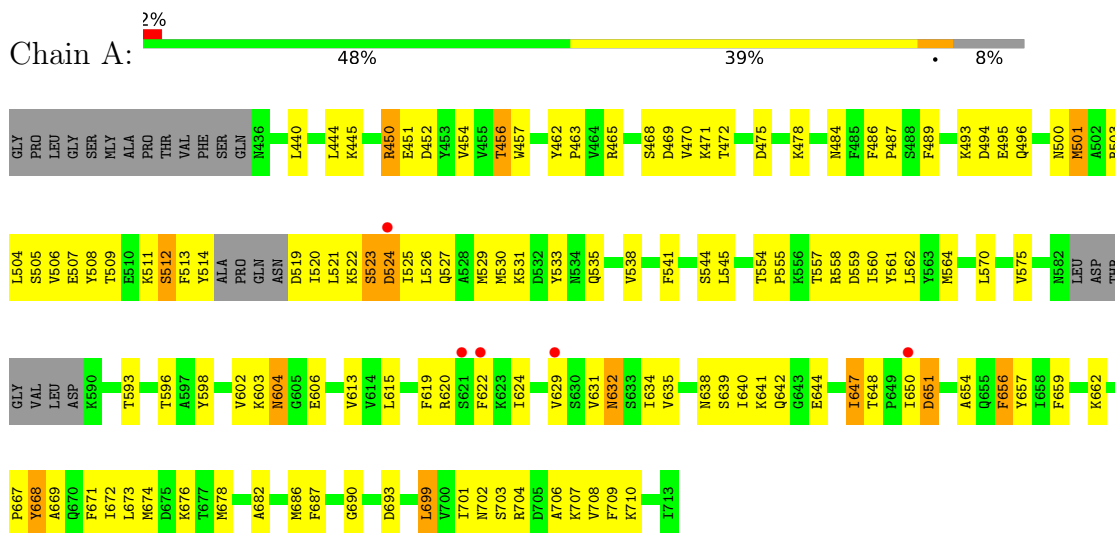
- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		

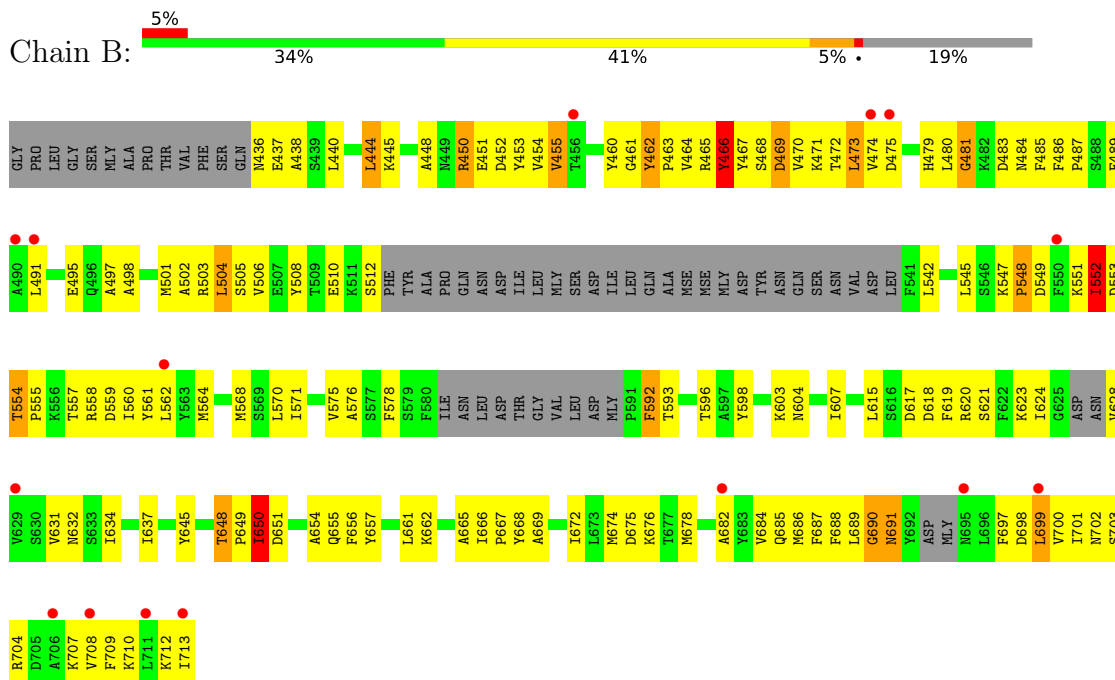
3 Residue-property plots [i](#)

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: General glycosylation pathway protein



- Molecule 1: General glycosylation pathway protein



4 Data and refinement statistics i

Property	Value	Source
Space group	P 64	Depositor
Cell constants a, b, c, α , β , γ	115.26Å 115.26Å 88.88Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.80 20.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (20.00-2.80) 99.6 (20.00-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.83 (at 2.80Å)	Xtrriage
Refinement program	REFMAC 5.5.0104	Depositor
R, R_{free}	0.245 , 0.260 0.244 , 0.260	Depositor DCC
R_{free} test set	1281 reflections (7.72%)	wwPDB-VP
Wilson B-factor (Å ²)	61.7	Xtrriage
Anisotropy	0.002	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 54.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.041 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4036	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% 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: CA, MLY

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	0.27	0/1997	0.57	0/2711
1	B	0.28	1/1753 (0.1%)	0.56	0/2379
All	All	0.27	1/3750 (0.0%)	0.56	0/5090

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	552	ILE	CB-CG1	-5.61	1.42	1.53

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	2176	0	2115	142	0
1	B	1859	0	1729	165	0
2	A	1	0	0	0	0
All	All	4036	0	3844	307	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 39.

The worst 5 of 307 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:504:LEU:HD23	1:B:504:LEU:O	1.45	1.15
1:A:501:MSE:HE3	1:A:501:MSE:HA	1.18	1.13
1:B:682:ALA:HB1	1:B:686:MSE:HE2	1.14	1.11
1:A:501:MSE:HA	1:A:501:MSE:CE	1.81	1.10
1:A:501:MSE:O	1:A:501:MSE:HE2	1.55	1.04

There are no symmetry-related clashes.

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	239/291 (82%)	209 (87%)	24 (10%)	6 (2%)	4	16
1	B	207/291 (71%)	159 (77%)	38 (18%)	10 (5%)	2	6
All	All	446/582 (77%)	368 (82%)	62 (14%)	16 (4%)	2	10

5 of 16 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	524	ASP
1	B	554	THR
1	B	466	TYR
1	B	552	ILE
1	B	690	GLY

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	215/228 (94%)	203 (94%)	12 (6%)	19	50
1	B	182/228 (80%)	168 (92%)	14 (8%)	12	36
All	All	397/456 (87%)	371 (94%)	26 (6%)	15	43

5 of 26 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	455	VAL
1	B	469	ASP
1	B	650	ILE
1	B	466	TYR
1	B	473	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	604	ASN
1	B	702	ASN
1	A	655	GLN
1	A	680	ASN
1	A	691	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

42 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	MLY	A	547	1	7,8,11	0.47	0	3,8,13	0.39	0
1	MLY	A	623	1	7,8,11	0.46	0	3,8,13	0.45	0
1	MLY	A	556	1	9,10,11	0.49	0	6,11,13	0.70	0
1	MLY	A	646	1	4,5,11	0.97	0	1,5,13	0.00	0
1	MLY	A	662	1	9,10,11	0.47	0	6,11,13	0.71	0
1	MLY	A	471	1	9,10,11	0.48	0	6,11,13	0.77	0
1	MLY	B	511	1	7,8,11	0.48	0	3,8,13	0.39	0
1	MLY	A	482	1	7,8,11	0.49	0	3,8,13	0.35	0
1	MLY	A	522	1	7,8,11	0.49	0	3,8,13	0.40	0
1	MLY	A	653	1	4,5,11	0.96	0	1,5,13	0.02	0
1	MLY	B	707	1	7,8,11	0.48	0	3,8,13	0.39	0
1	MLY	B	662	1	3,4,11	0.68	0	2,4,13	0.77	0
1	MLY	A	707	1	9,10,11	0.48	0	6,11,13	0.81	0
1	MLY	A	590	1	3,4,11	0.67	0	2,4,13	0.81	0
1	MLY	A	694	1	9,10,11	0.50	0	6,11,13	0.65	0
1	MLY	B	445	1	3,4,11	0.68	0	2,4,13	0.75	0
1	MLY	A	531	1	7,8,11	0.48	0	3,8,13	0.41	0
1	MLY	B	556	1	7,8,11	0.46	0	3,8,13	0.42	0
1	MLY	B	641	1	7,8,11	0.47	0	3,8,13	0.37	0
1	MLY	B	551	1	3,4,11	0.68	0	2,4,13	0.75	0
1	MLY	B	482	1	4,5,11	0.96	0	1,5,13	0.02	0
1	MLY	A	676	1	9,10,11	0.49	0	6,11,13	0.58	0
1	MLY	B	676	1	9,10,11	0.50	0	6,11,13	0.79	0
1	MLY	B	710	1	4,5,11	0.97	0	1,5,13	0.06	0
1	MLY	B	712	1	4,5,11	0.96	0	1,5,13	0.02	0
1	MLY	B	478	1	3,4,11	0.66	0	2,4,13	0.77	0
1	MLY	A	641	1	9,10,11	0.49	0	6,11,13	0.81	0
1	MLY	B	646	1	4,5,11	0.96	0	1,5,13	0.02	0
1	MLY	A	478	1	9,10,11	0.51	0	6,11,13	0.76	0
1	MLY	A	712	1	9,10,11	0.51	0	6,11,13	0.90	0
1	MLY	B	493	1	7,8,11	0.48	0	3,8,13	0.39	0
1	MLY	B	547	1	7,8,11	0.47	0	3,8,13	0.39	0
1	MLY	B	623	1	9,10,11	0.50	0	6,11,13	0.73	0
1	MLY	A	445	1	9,10,11	0.48	0	6,11,13	0.71	0
1	MLY	A	603	1	9,10,11	0.48	0	6,11,13	0.75	0
1	MLY	B	471	1	7,8,11	0.48	0	3,8,13	0.36	0
1	MLY	A	511	1	9,10,11	0.48	0	6,11,13	0.73	0
1	MLY	A	710	1	9,10,11	0.51	0	6,11,13	0.88	0
1	MLY	A	493	1	9,10,11	0.49	0	6,11,13	0.64	0
1	MLY	A	551	1	3,4,11	0.68	0	2,4,13	0.73	0
1	MLY	B	603	1	9,10,11	0.51	0	6,11,13	0.76	0
1	MLY	B	653	1	3,4,11	0.65	0	2,4,13	0.76	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	MLY	A	547	1	-	0/6/7/11	-
1	MLY	A	623	1	-	1/6/7/11	-
1	MLY	A	556	1	-	0/8/9/11	-
1	MLY	A	646	1	-	0/3/4/11	-
1	MLY	A	662	1	-	0/8/9/11	-
1	MLY	A	471	1	-	1/8/9/11	-
1	MLY	B	511	1	-	0/6/7/11	-
1	MLY	A	482	1	-	2/6/7/11	-
1	MLY	A	522	1	-	0/6/7/11	-
1	MLY	A	653	1	-	0/3/4/11	-
1	MLY	B	707	1	-	1/6/7/11	-
1	MLY	B	662	1	-	0/1/2/11	-
1	MLY	A	707	1	-	1/8/9/11	-
1	MLY	A	590	1	-	0/1/2/11	-
1	MLY	A	694	1	-	2/8/9/11	-
1	MLY	B	445	1	-	1/1/2/11	-
1	MLY	A	531	1	-	0/6/7/11	-
1	MLY	B	556	1	-	0/6/7/11	-
1	MLY	B	641	1	-	1/6/7/11	-
1	MLY	B	551	1	-	0/1/2/11	-
1	MLY	B	482	1	-	2/3/4/11	-
1	MLY	A	676	1	-	1/8/9/11	-
1	MLY	B	676	1	-	2/8/9/11	-
1	MLY	B	710	1	-	0/3/4/11	-
1	MLY	B	712	1	-	0/3/4/11	-
1	MLY	B	478	1	-	0/1/2/11	-
1	MLY	A	641	1	-	0/8/9/11	-
1	MLY	B	646	1	-	0/3/4/11	-
1	MLY	A	478	1	-	0/8/9/11	-
1	MLY	A	712	1	-	4/8/9/11	-
1	MLY	B	493	1	-	1/6/7/11	-
1	MLY	B	547	1	-	2/6/7/11	-
1	MLY	B	623	1	-	1/8/9/11	-
1	MLY	A	445	1	-	2/8/9/11	-
1	MLY	A	603	1	-	1/8/9/11	-
1	MLY	B	471	1	-	0/6/7/11	-
1	MLY	A	511	1	-	3/8/9/11	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	MLY	A	710	1	-	0/8/9/11	-
1	MLY	A	493	1	-	1/8/9/11	-
1	MLY	A	551	1	-	0/1/2/11	-
1	MLY	B	603	1	-	4/8/9/11	-
1	MLY	B	653	1	-	1/1/2/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 35 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	482	MLY	O-C-CA-CB
1	A	511	MLY	O-C-CA-CB
1	A	623	MLY	O-C-CA-CB
1	A	694	MLY	C-CA-CB-CG
1	A	707	MLY	O-C-CA-CB

There are no ring outliers.

24 monomers are involved in 47 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	662	MLY	1	0
1	A	471	MLY	2	0
1	A	522	MLY	2	0
1	B	707	MLY	3	0
1	B	662	MLY	1	0
1	A	707	MLY	7	0
1	B	445	MLY	1	0
1	A	531	MLY	2	0
1	B	551	MLY	1	0
1	A	676	MLY	1	0
1	B	676	MLY	2	0
1	B	710	MLY	3	0
1	B	712	MLY	2	0
1	A	641	MLY	1	0
1	A	478	MLY	2	0
1	B	547	MLY	4	0
1	B	623	MLY	1	0
1	A	445	MLY	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	603	MLY	1	0
1	B	471	MLY	1	0
1	A	511	MLY	3	0
1	A	710	MLY	1	0
1	A	493	MLY	2	0
1	B	603	MLY	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - 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 [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	236/291 (81%)	-0.05	5 (2%) 63 54	2, 20, 52, 59	0
1	B	211/291 (72%)	0.57	15 (7%) 22 16	6, 46, 71, 74	0
All	All	447/582 (76%)	0.25	20 (4%) 38 30	2, 35, 63, 74	0

The worst 5 of 20 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	474	VAL	4.0
1	B	695	ASN	3.8
1	B	711	LEU	3.1
1	A	621	SER	2.9
1	B	708	VAL	2.8

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	MLY	B	478	5/12	0.34	0.15	61,61,61,61	0
1	MLY	A	590	5/12	0.41	0.24	35,35,35,35	0
1	MLY	B	556	9/12	0.60	0.13	67,67,67,68	0
1	MLY	B	482	6/12	0.61	0.17	57,57,57,58	0
1	MLY	B	712	6/12	0.61	0.20	61,61,61,61	0
1	MLY	B	471	9/12	0.65	0.18	57,57,57,57	0
1	MLY	B	551	5/12	0.66	0.14	72,72,72,73	0
1	MLY	B	646	6/12	0.68	0.11	41,41,41,42	0
1	MLY	B	547	9/12	0.69	0.13	72,73,73,73	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
1	MLY	B	493	9/12	0.71	0.16	53,53,54,54	0
1	MLY	B	641	9/12	0.71	0.18	43,44,44,44	0
1	MLY	B	707	9/12	0.74	0.22	58,59,59,59	0
1	MLY	A	653	6/12	0.75	0.14	41,41,42,42	0
1	MLY	B	511	9/12	0.77	0.12	63,63,63,63	0
1	MLY	A	676	11/12	0.77	0.15	25,26,27,27	0
1	MLY	A	623	9/12	0.78	0.11	57,57,57,58	0
1	MLY	B	662	5/12	0.78	0.10	35,35,35,36	0
1	MLY	A	603	11/12	0.80	0.11	40,40,41,41	0
1	MLY	A	646	6/12	0.81	0.16	46,46,46,46	0
1	MLY	A	694	11/12	0.82	0.13	16,17,20,20	0
1	MLY	A	482	9/12	0.83	0.14	10,10,12,12	0
1	MLY	A	556	11/12	0.83	0.11	21,23,26,26	0
1	MLY	A	641	11/12	0.85	0.21	45,45,45,45	0
1	MLY	A	531	9/12	0.85	0.12	32,32,33,33	0
1	MLY	B	445	5/12	0.85	0.10	59,59,59,60	0
1	MLY	A	547	9/12	0.85	0.10	26,27,27,27	0
1	MLY	A	662	11/12	0.86	0.14	29,29,29,29	0
1	MLY	A	511	11/12	0.87	0.12	22,23,25,25	0
1	MLY	B	623	11/12	0.88	0.15	25,26,28,28	0
1	MLY	A	522	9/12	0.88	0.10	37,37,38,38	0
1	MLY	A	551	5/12	0.89	0.09	28,28,29,29	0
1	MLY	A	712	11/12	0.90	0.10	15,16,19,19	0
1	MLY	B	676	11/12	0.90	0.13	36,36,37,37	0
1	MLY	A	445	11/12	0.90	0.10	7,7,11,11	0
1	MLY	B	710	6/12	0.90	0.20	60,60,61,61	0
1	MLY	B	603	11/12	0.90	0.10	9,10,13,14	0
1	MLY	A	478	11/12	0.92	0.10	9,10,14,15	0
1	MLY	A	471	11/12	0.92	0.12	2,2,7,7	0
1	MLY	A	493	11/12	0.93	0.12	7,7,10,10	0
1	MLY	B	653	5/12	0.94	0.06	36,36,37,37	0
1	MLY	A	707	11/12	0.95	0.07	14,15,17,17	0
1	MLY	A	710	11/12	0.96	0.07	13,14,15,15	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	CA	A	1	1/1	0.95	0.04	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.