



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 01:43 PM UTC

PDB ID : 2VIX / pdb_00002vix
Title : Methylated Shigella flexneri MxiC
Authors : Deane, J.E.; Roversi, P.; King, C.; Johnson, S.; Lea, S.M.
Deposited on : 2007-12-05
Resolution : 2.85 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

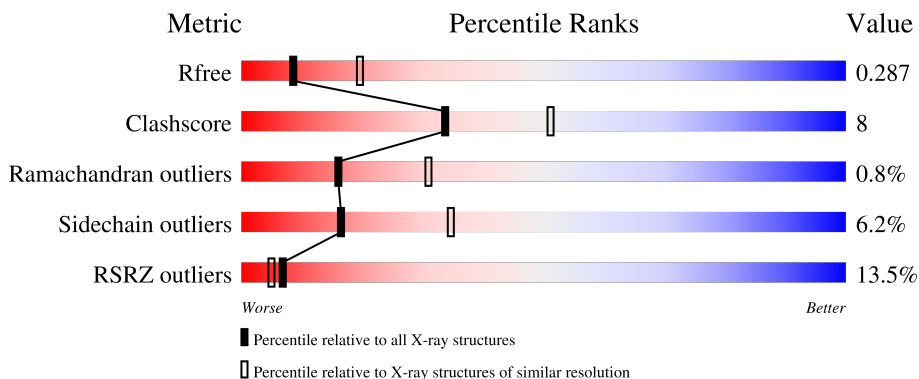
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	294	11% 81% 13% • •
1	B	294	5% 75% 19% • •
1	C	294	21% 72% 21% • •

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7117 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 PROTEIN MXIC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	283	Total	C	N	O	S	0	0	0
			2325	1489	381	447	8			
1	B	283	Total	C	N	O	S	0	0	0
			2325	1489	381	447	8			
1	C	281	Total	C	N	O	S	0	1	0
			2317	1487	378	444	8			

There are 36 discrepancies between the modelled and reference sequences:

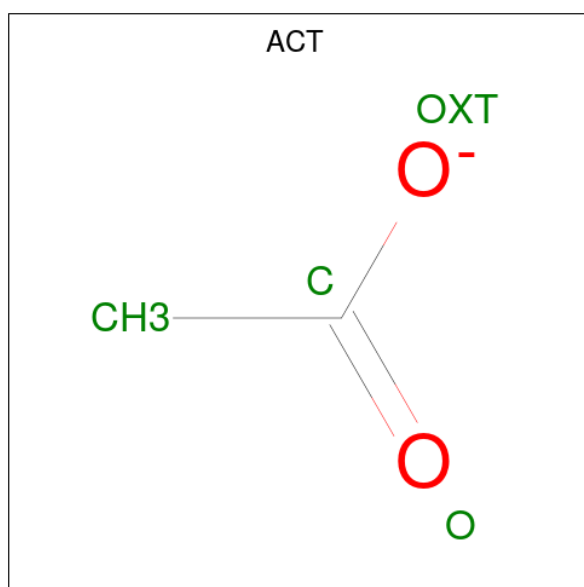
Chain	Residue	Modelled	Actual	Comment	Reference
A	62	HIS	-	expression tag	UNP Q04640
A	63	SER	-	expression tag	UNP Q04640
A	64	SER	-	expression tag	UNP Q04640
A	65	GLY	-	expression tag	UNP Q04640
A	66	LEU	-	expression tag	UNP Q04640
A	67	VAL	-	expression tag	UNP Q04640
A	68	PRO	-	expression tag	UNP Q04640
A	69	ARG	-	expression tag	UNP Q04640
A	70	GLY	-	expression tag	UNP Q04640
A	71	SER	-	expression tag	UNP Q04640
A	72	HIS	-	expression tag	UNP Q04640
A	73	MET	-	expression tag	UNP Q04640
B	62	HIS	-	expression tag	UNP Q04640
B	63	SER	-	expression tag	UNP Q04640
B	64	SER	-	expression tag	UNP Q04640
B	65	GLY	-	expression tag	UNP Q04640
B	66	LEU	-	expression tag	UNP Q04640
B	67	VAL	-	expression tag	UNP Q04640
B	68	PRO	-	expression tag	UNP Q04640
B	69	ARG	-	expression tag	UNP Q04640
B	70	GLY	-	expression tag	UNP Q04640
B	71	SER	-	expression tag	UNP Q04640
B	72	HIS	-	expression tag	UNP Q04640

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Chain	Residue	Modelled	Actual	Comment	Reference
B	73	MET	-	expression tag	UNP Q04640
C	62	HIS	-	expression tag	UNP Q04640
C	63	SER	-	expression tag	UNP Q04640
C	64	SER	-	expression tag	UNP Q04640
C	65	GLY	-	expression tag	UNP Q04640
C	66	LEU	-	expression tag	UNP Q04640
C	67	VAL	-	expression tag	UNP Q04640
C	68	PRO	-	expression tag	UNP Q04640
C	69	ARG	-	expression tag	UNP Q04640
C	70	GLY	-	expression tag	UNP Q04640
C	71	SER	-	expression tag	UNP Q04640
C	72	HIS	-	expression tag	UNP Q04640
C	73	MET	-	expression tag	UNP Q04640

- Molecule 2 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

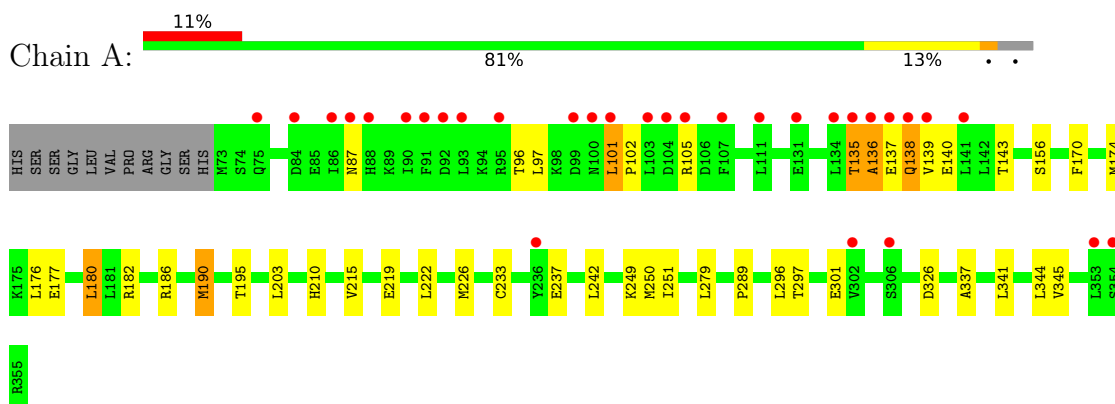
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	46	Total	O	0	0
			46	46		
4	B	46	Total	O	0	0
			46	46		
4	C	42	Total	O	0	0
			42	42		

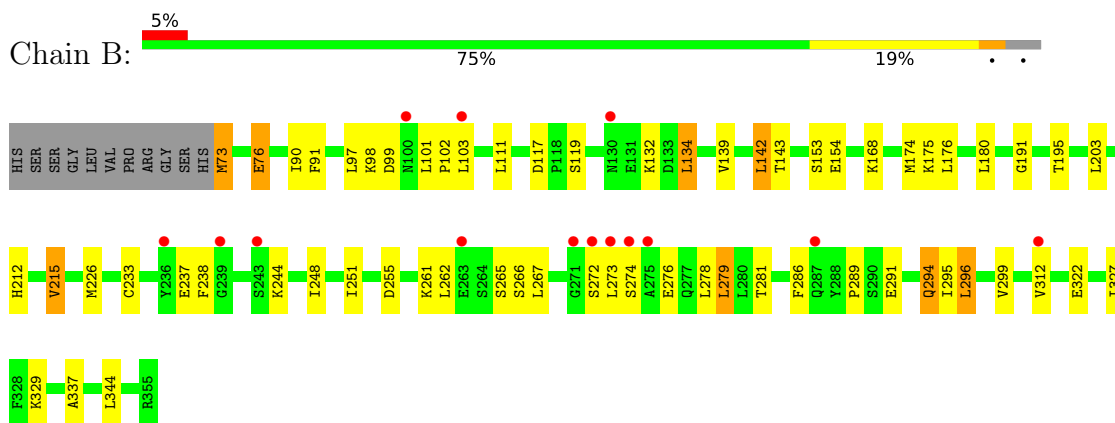
3 Residue-property plots [i](#)

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.

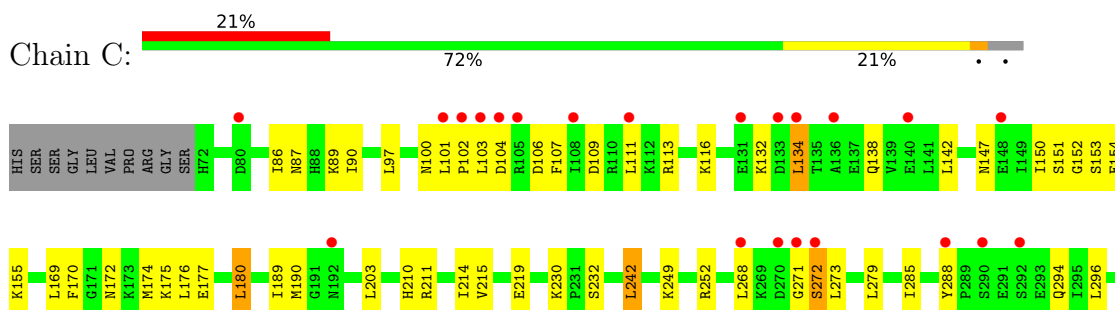
• Molecule 1: PROTEIN MXIC

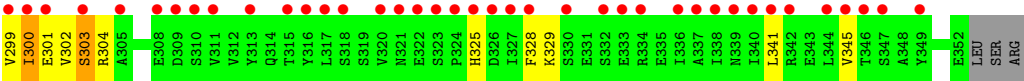


• Molecule 1: PROTEIN MXIC



• Molecule 1: PROTEIN MXIC





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.79Å 102.57Å 122.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.40 – 2.85 50.40 – 2.85	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.40-2.85) 91.8 (50.40-2.85)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.86Å)	Xtriage
Refinement program	TNT 5.13.1.0	Depositor
R, R_{free}	0.244 , 0.273 0.255 , 0.287	Depositor DCC
R_{free} test set	1216 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	52.5	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	7117	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.35% 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: GOL, ACT, 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	0.30	0/2139	0.67	1/2901 (0.0%)
1	B	0.30	0/2139	0.66	0/2901
1	C	0.30	0/2137	0.68	1/2901 (0.0%)
All	All	0.30	0/6415	0.67	2/8703 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	138	GLN	N-CA-C	-6.22	105.67	113.20
1	C	106	ASP	N-CA-C	-5.76	106.25	113.28

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	2325	0	2362	23	0
1	B	2325	0	2363	41	0
1	C	2317	0	2346	46	0
2	A	4	0	3	0	0
3	A	6	0	8	0	0
3	B	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	46	0	0	0	0
4	B	46	0	0	0	0
4	C	42	0	0	0	0
All	All	7117	0	7090	109	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 8.

The worst 5 of 109 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:132:MLY:HA	1:B:134:LEU:H	1.10	1.11
1:B:102:PRO:HA	1:B:103:LEU:HB2	1.29	1.08
1:B:132:MLY:HA	1:B:134:LEU:N	1.83	0.93
1:B:102:PRO:HA	1:B:103:LEU:CB	2.07	0.85
1:C:328:PHE:N	1:C:329:MLY:HA	1.92	0.84

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	261/294 (89%)	249 (95%)	9 (3%)	3 (1%)	11	24
1	B	261/294 (89%)	252 (97%)	9 (3%)	0	100	100
1	C	260/294 (88%)	242 (93%)	15 (6%)	3 (1%)	10	22
All	All	782/882 (89%)	743 (95%)	33 (4%)	6 (1%)	16	31

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	137	GLU
1	A	97	LEU
1	C	271	GLY
1	A	136	ALA
1	C	303	SER

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	241/250 (96%)	227 (94%)	14 (6%)	18	38
1	B	241/250 (96%)	224 (93%)	17 (7%)	13	28
1	C	240/250 (96%)	226 (94%)	14 (6%)	18	38
All	All	722/750 (96%)	677 (94%)	45 (6%)	16	34

5 of 45 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	296	LEU
1	C	180	LEU
1	B	312	VAL
1	C	134	LEU
1	C	242	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	210	HIS
1	C	339	ASN
1	B	166	GLN
1	B	287	GLN
1	C	100	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

60 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	144	1	9,10,11	0.52	0	6,11,13	0.84	0
1	MLY	B	329	1	9,10,11	0.52	0	6,11,13	0.78	0
1	MLY	C	329	1	9,10,11	0.56	0	6,11,13	0.70	0
1	MLY	A	173	1	9,10,11	0.53	0	6,11,13	0.78	0
1	MLY	B	244	1	9,10,11	0.57	0	6,11,13	0.74	0
1	MLY	A	175	1	9,10,11	0.59	0	6,11,13	0.69	0
1	MLY	C	144	1	9,10,11	0.51	0	6,11,13	0.81	0
1	MLY	A	230	1	9,10,11	0.56	0	6,11,13	0.72	0
1	MLY	B	112	1	9,10,11	0.53	0	6,11,13	0.86	0
1	MLY	B	144	1	9,10,11	0.50	0	6,11,13	0.73	0
1	MLY	A	155	1	9,10,11	0.52	0	6,11,13	0.84	0
1	MLY	B	269	1	9,10,11	0.55	0	6,11,13	0.76	0
1	MLY	B	350	1	9,10,11	0.47	0	6,11,13	0.79	0
1	MLY	C	98	1	9,10,11	0.55	0	6,11,13	0.75	0
1	MLY	C	116	1	9,10,11	0.56	0	6,11,13	0.72	0
1	MLY	C	132	1	9,10,11	0.45	0	6,11,13	0.90	0
1	MLY	B	116	1	9,10,11	0.57	0	6,11,13	0.75	0
1	MLY	B	173	1	9,10,11	0.54	0	6,11,13	0.81	0
1	MLY	C	89	1	9,10,11	0.55	0	6,11,13	0.72	0
1	MLY	B	132	1	9,10,11	0.53	0	6,11,13	0.77	0
1	MLY	A	351	1	9,10,11	0.60	0	6,11,13	0.74	0
1	MLY	A	329	1	9,10,11	0.57	0	6,11,13	0.71	0
1	MLY	B	249	1	9,10,11	0.59	0	6,11,13	0.78	0
1	MLY	C	351	1	9,10,11	0.57	0	6,11,13	0.71	0
1	MLY	B	94	1	9,10,11	0.55	0	6,11,13	0.77	0
1	MLY	C	230	1	9,10,11	0.54	0	6,11,13	0.77	0
1	MLY	A	112	1	9,10,11	0.58	0	6,11,13	0.75	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	MLY	A	269	1	9,10,11	0.51	0	6,11,13	0.72	0
1	MLY	C	155	1	9,10,11	0.50	0	6,11,13	0.75	0
1	MLY	C	261	1	9,10,11	0.65	0	6,11,13	0.73	0
1	MLY	B	175	1	9,10,11	0.56	0	6,11,13	0.71	0
1	MLY	A	89	1	9,10,11	0.61	0	6,11,13	0.75	0
1	MLY	B	351	1	9,10,11	0.52	0	6,11,13	0.73	0
1	MLY	B	230	1	9,10,11	0.56	0	6,11,13	0.69	0
1	MLY	A	260	1	9,10,11	0.47	0	6,11,13	0.78	0
1	MLY	B	155	1	9,10,11	0.54	0	6,11,13	0.72	0
1	MLY	A	116	1	9,10,11	0.56	0	6,11,13	0.74	0
1	MLY	C	112	1	9,10,11	0.58	0	6,11,13	0.72	0
1	MLY	C	269	1	9,10,11	0.61	0	6,11,13	0.72	0
1	MLY	A	261	1	9,10,11	0.51	0	6,11,13	0.96	0
1	MLY	B	260	1	9,10,11	0.55	0	6,11,13	0.79	0
1	MLY	C	244	1	9,10,11	0.59	0	6,11,13	0.77	0
1	MLY	C	260	1	9,10,11	0.56	0	6,11,13	0.82	0
1	MLY	B	168	1	9,10,11	0.53	0	6,11,13	0.78	0
1	MLY	C	94	1	9,10,11	0.54	0	6,11,13	0.72	0
1	MLY	C	249	1	9,10,11	0.56	0	6,11,13	0.78	0
1	MLY	A	244	1	9,10,11	0.53	0	6,11,13	0.94	0
1	MLY	B	89	1	9,10,11	0.60	0	6,11,13	0.72	0
1	MLY	B	98	1	9,10,11	0.54	0	6,11,13	0.88	0
1	MLY	B	261	1	9,10,11	0.62	0	6,11,13	0.77	0
1	MLY	C	168	1	9,10,11	0.53	0	6,11,13	0.81	0
1	MLY	C	173	1	9,10,11	0.59	0	6,11,13	0.74	0
1	MLY	A	94	1	9,10,11	0.57	0	6,11,13	0.76	0
1	MLY	A	350	1	9,10,11	0.58	0	6,11,13	0.73	0
1	MLY	C	175	1	9,10,11	0.57	0	6,11,13	0.81	0
1	MLY	A	98	1	9,10,11	0.54	0	6,11,13	0.77	0
1	MLY	A	168	1	9,10,11	0.56	0	6,11,13	0.78	0
1	MLY	C	350	1	9,10,11	0.49	0	6,11,13	0.73	0
1	MLY	A	132	1	9,10,11	0.51	0	6,11,13	0.91	0
1	MLY	A	249	1	9,10,11	0.63	0	6,11,13	0.72	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	144	1	-	3/8/9/11	-
1	MLY	B	329	1	-	3/8/9/11	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	MLY	C	329	1	-	6/8/9/11	-
1	MLY	A	173	1	-	6/8/9/11	-
1	MLY	B	244	1	-	4/8/9/11	-
1	MLY	A	175	1	-	3/8/9/11	-
1	MLY	C	144	1	-	4/8/9/11	-
1	MLY	A	230	1	-	3/8/9/11	-
1	MLY	B	112	1	-	2/8/9/11	-
1	MLY	B	144	1	-	3/8/9/11	-
1	MLY	A	155	1	-	4/8/9/11	-
1	MLY	B	269	1	-	5/8/9/11	-
1	MLY	B	350	1	-	0/8/9/11	-
1	MLY	C	98	1	-	5/8/9/11	-
1	MLY	C	116	1	-	8/8/9/11	-
1	MLY	C	132	1	-	4/8/9/11	-
1	MLY	B	116	1	-	4/8/9/11	-
1	MLY	B	173	1	-	3/8/9/11	-
1	MLY	C	89	1	-	2/8/9/11	-
1	MLY	B	132	1	-	6/8/9/11	-
1	MLY	A	351	1	-	6/8/9/11	-
1	MLY	A	329	1	-	4/8/9/11	-
1	MLY	B	249	1	-	3/8/9/11	-
1	MLY	C	351	1	-	7/8/9/11	-
1	MLY	B	94	1	-	4/8/9/11	-
1	MLY	C	230	1	-	4/8/9/11	-
1	MLY	A	112	1	-	6/8/9/11	-
1	MLY	A	269	1	-	3/8/9/11	-
1	MLY	C	155	1	-	6/8/9/11	-
1	MLY	C	261	1	-	3/8/9/11	-
1	MLY	B	175	1	-	5/8/9/11	-
1	MLY	A	89	1	-	3/8/9/11	-
1	MLY	B	351	1	-	6/8/9/11	-
1	MLY	B	230	1	-	5/8/9/11	-
1	MLY	A	260	1	-	1/8/9/11	-
1	MLY	B	155	1	-	4/8/9/11	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	MLY	A	116	1	-	6/8/9/11	-
1	MLY	C	112	1	-	3/8/9/11	-
1	MLY	C	269	1	-	6/8/9/11	-
1	MLY	A	261	1	-	1/8/9/11	-
1	MLY	B	260	1	-	4/8/9/11	-
1	MLY	C	244	1	-	4/8/9/11	-
1	MLY	C	260	1	-	3/8/9/11	-
1	MLY	B	168	1	-	3/8/9/11	-
1	MLY	C	94	1	-	5/8/9/11	-
1	MLY	C	249	1	-	5/8/9/11	-
1	MLY	A	244	1	-	2/8/9/11	-
1	MLY	B	89	1	-	0/8/9/11	-
1	MLY	B	98	1	-	2/8/9/11	-
1	MLY	B	261	1	-	3/8/9/11	-
1	MLY	C	168	1	-	5/8/9/11	-
1	MLY	C	173	1	-	3/8/9/11	-
1	MLY	A	94	1	-	3/8/9/11	-
1	MLY	A	350	1	-	5/8/9/11	-
1	MLY	C	175	1	-	2/8/9/11	-
1	MLY	A	98	1	-	4/8/9/11	-
1	MLY	A	168	1	-	3/8/9/11	-
1	MLY	C	350	1	-	3/8/9/11	-
1	MLY	A	132	1	-	4/8/9/11	-
1	MLY	A	249	1	-	4/8/9/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 231 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	112	MLY	O-C-CA-CB
1	A	116	MLY	N-CA-CB-CG
1	A	116	MLY	C-CA-CB-CG
1	A	132	MLY	N-CA-CB-CG
1	A	132	MLY	C-CA-CB-CG

There are no ring outliers.

16 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	329	MLY	1	0
1	C	329	MLY	3	0
1	B	244	MLY	1	0
1	C	116	MLY	1	0
1	C	132	MLY	1	0
1	C	89	MLY	1	0
1	B	132	MLY	4	0
1	C	230	MLY	1	0
1	C	155	MLY	1	0
1	B	175	MLY	1	0
1	B	168	MLY	1	0
1	C	249	MLY	1	0
1	B	98	MLY	2	0
1	B	261	MLY	1	0
1	C	175	MLY	2	0
1	A	249	MLY	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	ACT	A	1356	-	3,3,3	1.21	0	3,3,3	0.90	0
3	GOL	B	1356	-	5,5,5	0.48	0	5,5,5	1.79	2 (40%)
3	GOL	A	1357	-	5,5,5	0.32	0	5,5,5	1.70	1 (20%)

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
3	GOL	B	1356	-	-	1/4/4/4	-
3	GOL	A	1357	-	-	2/4/4/4	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	B	1356	GOL	O1-C1-C2	-2.71	98.16	110.38
3	A	1357	GOL	O3-C3-C2	-2.36	99.76	110.38
3	B	1356	GOL	C3-C2-C1	-2.21	103.71	111.80

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1357	GOL	C1-C2-C3-O3
3	B	1356	GOL	O1-C1-C2-C3
3	A	1357	GOL	O2-C2-C3-O3

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	263/294 (89%)	0.79	31 (11%) 9 7	32, 53, 96, 103	0
1	B	263/294 (89%)	0.68	14 (5%) 32 24	31, 50, 70, 83	0
1	C	261/294 (88%)	1.22	61 (23%) 2 2	21, 65, 121, 131	1 (0%)
All	All	787/882 (89%)	0.90	106 (13%) 7 5	21, 53, 112, 131	1 (0%)

The worst 5 of 106 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	136	ALA	5.5
1	A	93	LEU	4.9
1	C	327	ILE	4.9
1	C	134	LEU	4.7
1	C	345	VAL	4.7

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	C	329	11/12	0.55	0.28	112,115,124,130	0
1	MLY	C	350	11/12	0.55	0.29	119,124,125,126	0
1	MLY	B	269	11/12	0.57	0.24	73,78,83,83	0
1	MLY	A	132	11/12	0.58	0.23	81,92,106,107	0
1	MLY	A	98	11/12	0.58	0.22	83,96,102,106	0
1	MLY	C	351	11/12	0.62	0.22	121,124,126,130	0
1	MLY	C	269	11/12	0.63	0.29	81,85,93,99	0
1	MLY	A	116	11/12	0.64	0.23	60,81,99,104	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	MLY	A	144	11/12	0.65	0.19	75,78,93,93	0
1	MLY	C	98	11/12	0.65	0.27	80,85,97,98	0
1	MLY	B	98	11/12	0.66	0.25	56,66,77,79	0
1	MLY	A	269	11/12	0.67	0.24	64,68,87,89	0
1	MLY	C	144	11/12	0.67	0.23	64,72,85,85	0
1	MLY	C	132	11/12	0.68	0.22	75,85,93,104	0
1	MLY	A	351	11/12	0.69	0.22	62,73,87,87	0
1	MLY	B	132	11/12	0.73	0.28	54,62,84,98	0
1	MLY	B	351	11/12	0.76	0.22	49,62,81,90	0
1	MLY	A	350	11/12	0.76	0.20	63,76,82,85	0
1	MLY	C	116	11/12	0.76	0.20	49,52,83,83	0
1	MLY	C	112	11/12	0.77	0.23	51,68,80,91	0
1	MLY	C	230	11/12	0.79	0.20	39,49,74,82	0
1	MLY	C	94	11/12	0.81	0.20	53,69,75,76	0
1	MLY	C	261	11/12	0.81	0.15	59,64,76,80	0
1	MLY	B	116	11/12	0.82	0.15	30,39,66,73	0
1	MLY	A	112	11/12	0.82	0.21	70,72,83,92	0
1	MLY	B	261	11/12	0.83	0.17	53,58,70,81	0
1	MLY	A	94	11/12	0.83	0.22	77,92,98,98	0
1	MLY	A	261	11/12	0.83	0.15	36,49,67,67	0
1	MLY	B	168	11/12	0.83	0.16	37,46,57,59	0
1	MLY	B	175	11/12	0.83	0.15	44,52,79,79	0
1	MLY	B	144	11/12	0.84	0.15	35,42,64,64	0
1	MLY	A	89	11/12	0.85	0.20	64,80,91,92	0
1	MLY	B	173	11/12	0.85	0.14	23,54,58,58	0
1	MLY	B	329	11/12	0.85	0.14	48,58,79,81	0
1	MLY	B	155	11/12	0.85	0.21	40,43,57,57	0
1	MLY	B	260	11/12	0.86	0.17	48,55,73,80	0
1	MLY	C	168	11/12	0.86	0.14	20,37,50,64	0
1	MLY	A	329	11/12	0.86	0.14	48,59,79,79	0
1	MLY	B	350	11/12	0.86	0.15	14,46,65,65	0
1	MLY	A	173	11/12	0.87	0.17	38,45,63,69	0
1	MLY	C	260	11/12	0.88	0.14	40,56,65,71	0
1	MLY	B	112	11/12	0.88	0.16	31,43,58,58	0
1	MLY	A	230	11/12	0.88	0.15	18,48,76,85	0
1	MLY	A	168	11/12	0.90	0.15	18,34,56,59	0
1	MLY	C	249	11/12	0.91	0.12	14,36,42,54	0
1	MLY	B	249	11/12	0.91	0.12	24,45,50,68	0
1	MLY	A	260	11/12	0.91	0.14	29,43,74,75	0
1	MLY	A	155	11/12	0.91	0.12	34,44,51,53	0
1	MLY	B	94	11/12	0.91	0.15	50,52,67,69	0
1	MLY	B	230	11/12	0.91	0.14	34,47,70,71	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	MLY	C	244	11/12	0.91	0.14	33,40,76,76	0
1	MLY	A	175	11/12	0.92	0.11	15,40,53,53	0
1	MLY	C	173	11/12	0.92	0.14	26,39,63,66	0
1	MLY	C	175	11/12	0.92	0.11	20,35,42,42	0
1	MLY	B	244	11/12	0.92	0.12	33,41,45,45	0
1	MLY	A	249	11/12	0.92	0.13	16,35,46,46	0
1	MLY	C	155	11/12	0.92	0.15	44,47,62,62	0
1	MLY	A	244	11/12	0.93	0.13	21,33,62,62	0
1	MLY	B	89	11/12	0.94	0.11	14,41,52,59	0
1	MLY	C	89	11/12	0.95	0.13	41,62,63,66	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($\text{\AA}^2$)	Q<0.9
3	GOL	A	1357	6/6	0.40	0.25	87,87,87,87	0
2	ACT	A	1356	4/4	0.61	0.32	95,95,95,95	0
3	GOL	B	1356	6/6	0.70	0.27	90,90,90,90	0

6.5 Other polymers [i](#)

There are no such residues in this entry.