



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 03:18 AM UTC

PDB ID : 7JVH / pdb_00007jvh
Title : Crystal structure of a GH43_12 retrieved from capybara gut metagenome
Authors : Cabral, L.; Domingues, M.N.; Martins, M.P.; Persinoti, G.F.; Morais, M.A.B.;
Murakami, M.T.
Deposited on : 2020-08-21
Resolution : 2.48 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

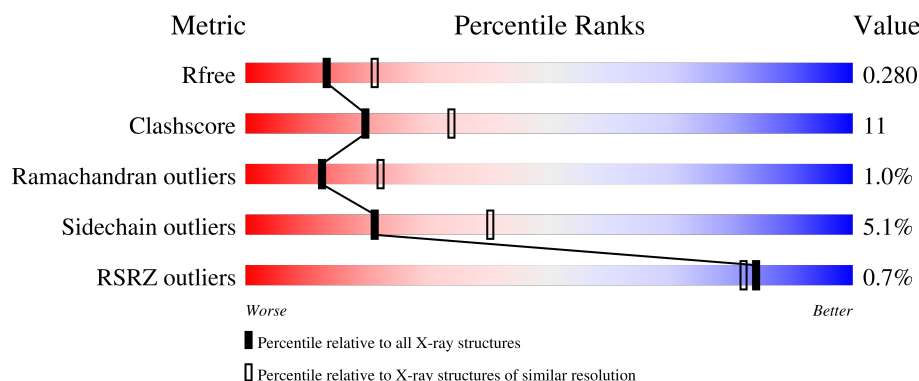
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.48 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	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (2.50-2.46)

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	567	
1	B	567	
1	C	567	
1	D	567	
1	E	567	

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Mol	Chain	Length	Quality of chain
1	F	567	2% 63% 20% • 12%

2 Entry composition [i](#)

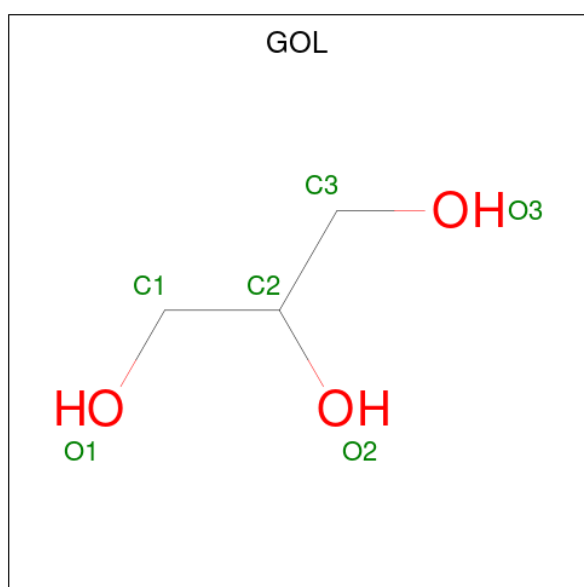
There are 3 unique types of molecules in this entry. The entry contains 25078 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glycoside Hydrolase Family 43_12.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	535	Total	C	N	O	S	Se	0	0	0
			4232	2721	700	793	9	9			
1	B	535	Total	C	N	O	S	Se	0	0	0
			4231	2719	700	794	9	9			
1	C	536	Total	C	N	O	S	Se	0	0	0
			4239	2725	702	794	9	9			
1	D	534	Total	C	N	O	S	Se	0	0	0
			4221	2713	698	792	9	9			
1	E	497	Total	C	N	O	S	Se	0	0	0
			3943	2543	650	733	8	9			
1	F	499	Total	C	N	O	S	Se	0	0	0
			3964	2556	651	740	8	9			

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

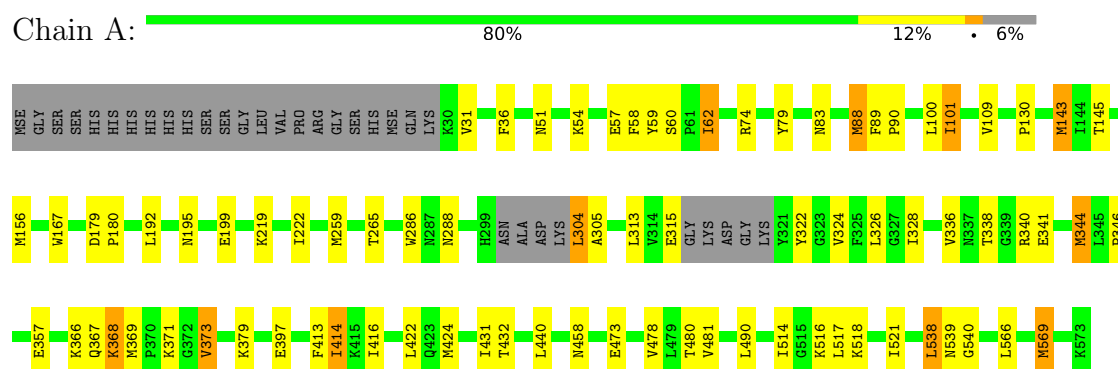
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	64	Total O 64 64	0	0
3	B	52	Total O 52 52	0	0
3	C	32	Total O 32 32	0	0
3	D	31	Total O 31 31	0	0
3	E	4	Total O 4 4	0	0
3	F	5	Total O 5 5	0	0

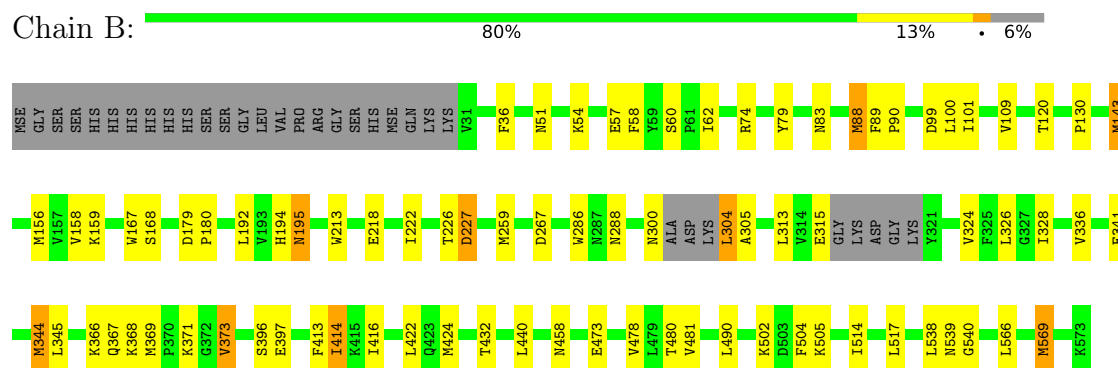
3 Residue-property plots [i](#)

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.

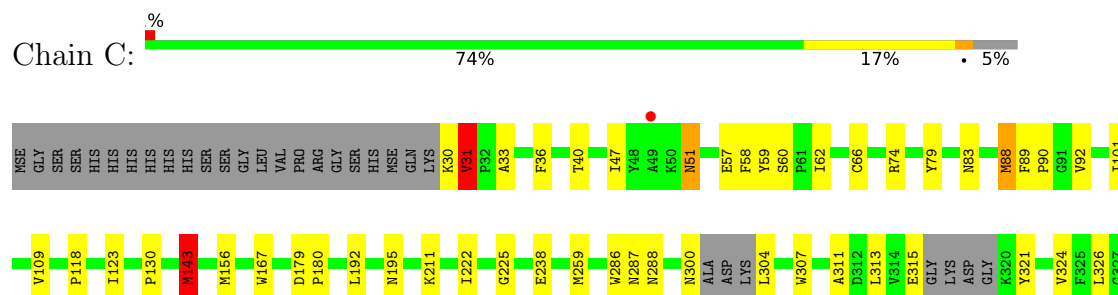
• Molecule 1: Glycoside Hydrolase Family 43_12

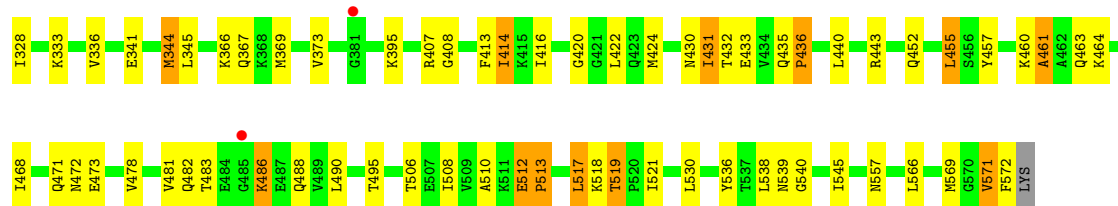


• Molecule 1: Glycoside Hydrolase Family 43_12



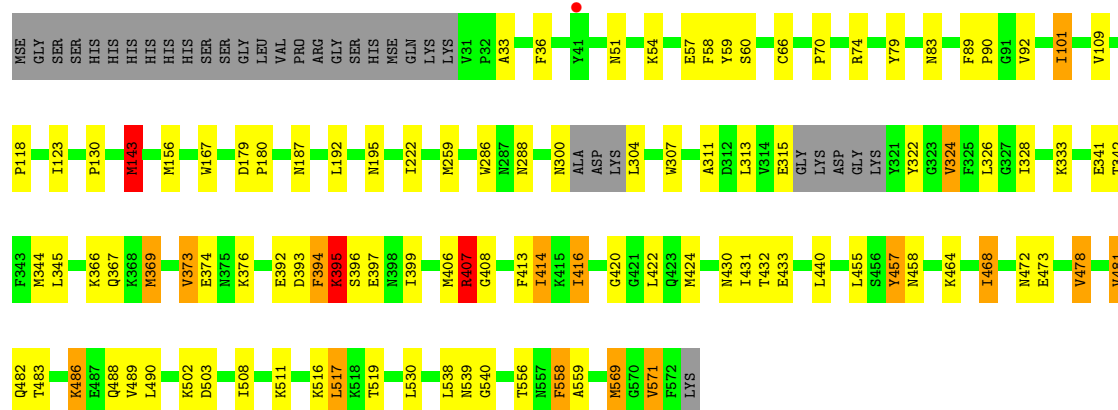
• Molecule 1: Glycoside Hydrolase Family 43_12





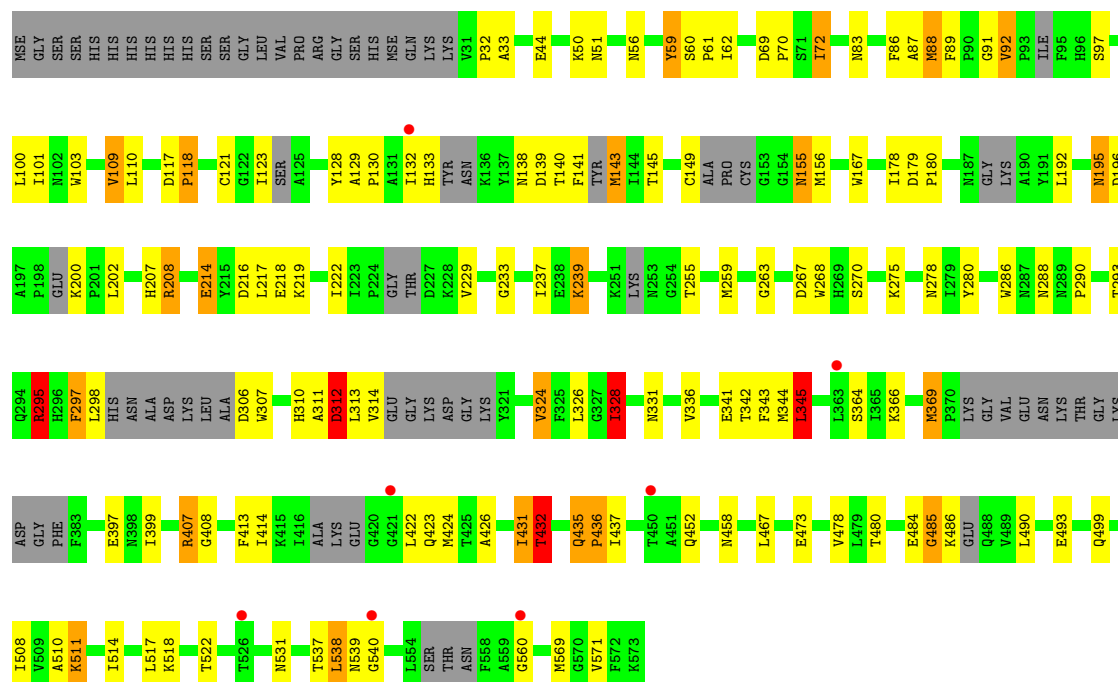
• Molecule 1: Glycoside Hydrolase Family 43_12

Chain D: 75% 15% 6%

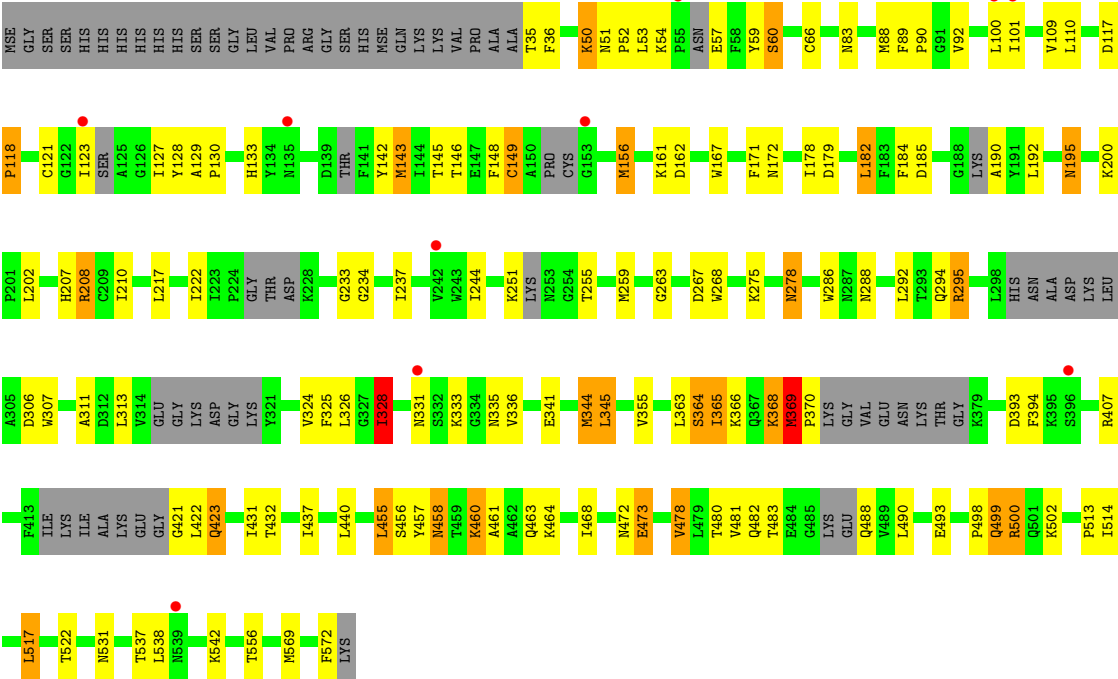


• Molecule 1: Glycoside Hydrolase Family 43_12

Chain E: 62% 20% 12%



• Molecule 1: Glycoside Hydrolase Family 43_12



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	107.22Å 107.38Å 132.53Å 99.03° 113.38° 108.45°	Depositor
Resolution (Å)	49.20 – 2.48 49.20 – 2.48	Depositor EDS
% Data completeness (in resolution range)	96.0 (49.20-2.48) 95.9 (49.20-2.48)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 2.48Å)	Xtriage
Refinement program	PHENIX 1.8	Depositor
R, R_{free}	0.246 , 0.278 0.250 , 0.280	Depositor DCC
R_{free} test set	8303 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	53.8	Xtriage
Anisotropy	0.711	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.437 for k,h,-h-k-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	25078	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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: GOL

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	0.65	1/4343 (0.0%)	0.87	3/5887 (0.1%)
1	B	0.67	2/4342 (0.0%)	0.91	6/5887 (0.1%)
1	C	0.80	7/4350 (0.2%)	1.10	21/5898 (0.4%)
1	D	0.74	4/4332 (0.1%)	1.07	19/5876 (0.3%)
1	E	0.79	6/4037 (0.1%)	1.12	27/5459 (0.5%)
1	F	0.73	3/4062 (0.1%)	1.10	20/5498 (0.4%)
All	All	0.73	23/25466 (0.1%)	1.03	96/34505 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2

The worst 5 of 23 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	513	PRO	N-CA	18.04	1.69	1.47
1	F	118	PRO	N-CA	17.89	1.70	1.47
1	E	436	PRO	N-CA	17.48	1.69	1.47
1	E	118	PRO	N-CA	15.56	1.67	1.47
1	F	345	LEU	C-N	13.52	1.51	1.33

The worst 5 of 96 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	512	GLU	CA-C-N	17.45	138.22	119.90
1	C	512	GLU	C-N-CA	17.45	138.22	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	161	LYS	N-CA-C	-14.87	84.49	108.73
1	F	117	ASP	CA-C-N	14.59	138.07	119.84
1	F	117	ASP	C-N-CA	14.59	138.07	119.84

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	226	THR	Mainchain
1	B	227	ASP	Mainchain

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	4232	0	4103	54	0
1	B	4231	0	4096	57	0
1	C	4239	0	4109	97	0
1	D	4221	0	4083	87	0
1	E	3943	0	3807	135	0
1	F	3964	0	3804	140	0
2	A	18	0	24	1	0
2	B	12	0	16	0	0
2	C	12	0	16	3	0
2	D	12	0	16	0	0
2	F	6	0	8	0	0
3	A	64	0	0	0	0
3	B	52	0	0	0	0
3	C	32	0	0	0	0
3	D	31	0	0	0	0
3	E	4	0	0	0	0
3	F	5	0	0	0	0
All	All	25078	0	24082	562	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 11.

The worst 5 of 562 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:E:436:PRO:N	1:E:436:PRO:CA	1.69	1.42
1:C:513:PRO:N	1:C:513:PRO:CA	1.69	1.38
1:F:118:PRO:N	1:F:118:PRO:CA	1.70	1.34
1:F:36:PHE:CE1	1:F:455:LEU:HD13	1.69	1.25
1:E:118:PRO:HB2	1:E:407:ARG:HA	1.15	1.11

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	529/567 (93%)	498 (94%)	28 (5%)	3 (1%)	21	35
1	B	529/567 (93%)	497 (94%)	30 (6%)	2 (0%)	30	46
1	C	530/567 (94%)	500 (94%)	26 (5%)	4 (1%)	16	28
1	D	528/567 (93%)	489 (93%)	31 (6%)	8 (2%)	8	14
1	E	465/567 (82%)	425 (91%)	34 (7%)	6 (1%)	9	16
1	F	473/567 (83%)	438 (93%)	28 (6%)	7 (2%)	8	14
All	All	3054/3402 (90%)	2847 (93%)	177 (6%)	30 (1%)	12	22

5 of 30 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	31	VAL
1	A	539	ASN
1	B	539	ASN
1	C	539	ASN
1	D	395	LYS

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	455/470 (97%)	439 (96%)	16 (4%)	32	55
1	B	455/470 (97%)	442 (97%)	13 (3%)	37	62
1	C	456/470 (97%)	439 (96%)	17 (4%)	30	54
1	D	454/470 (97%)	430 (95%)	24 (5%)	20	39
1	E	425/470 (90%)	388 (91%)	37 (9%)	9	18
1	F	426/470 (91%)	397 (93%)	29 (7%)	14	28
All	All	2671/2820 (95%)	2535 (95%)	136 (5%)	21	40

5 of 136 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	328	ILE
1	F	369	MSE
1	F	500	ARG
1	D	156	MSE
1	D	143	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 21 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	488	GLN
1	F	248	HIS
1	F	539	ASN
1	F	458	ASN
1	E	488	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

10 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	GOL	B	601	-	5,5,5	0.12	0	5,5,5	0.40	0
2	GOL	D	602	-	5,5,5	0.11	0	5,5,5	0.37	0
2	GOL	C	601	-	5,5,5	0.11	0	5,5,5	0.34	0
2	GOL	F	601	-	5,5,5	0.10	0	5,5,5	0.26	0
2	GOL	B	602	-	5,5,5	0.12	0	5,5,5	0.36	0
2	GOL	C	602	-	5,5,5	0.16	0	5,5,5	0.37	0
2	GOL	A	601	-	5,5,5	0.10	0	5,5,5	0.27	0
2	GOL	A	602	-	5,5,5	0.09	0	5,5,5	0.29	0
2	GOL	D	601	-	5,5,5	0.10	0	5,5,5	0.33	0
2	GOL	A	603	-	5,5,5	0.09	0	5,5,5	0.24	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
2	GOL	B	601	-	-	0/4/4/4	-
2	GOL	D	602	-	-	3/4/4/4	-
2	GOL	C	601	-	-	2/4/4/4	-
2	GOL	F	601	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	B	602	-	-	2/4/4/4	-
2	GOL	C	602	-	-	3/4/4/4	-
2	GOL	A	601	-	-	1/4/4/4	-
2	GOL	A	602	-	-	2/4/4/4	-
2	GOL	D	601	-	-	2/4/4/4	-
2	GOL	A	603	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 18 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	601	GOL	C1-C2-C3-O3
2	D	601	GOL	O1-C1-C2-C3
2	D	601	GOL	O1-C1-C2-O2
2	A	602	GOL	C1-C2-C3-O3
2	A	603	GOL	O1-C1-C2-C3

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	602	GOL	3	0
2	A	602	GOL	1	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 [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	526/567 (92%)	-0.60	0 100 100	41, 58, 83, 112	0
1	B	526/567 (92%)	-0.64	0 100 100	41, 58, 85, 115	0
1	C	527/567 (92%)	-0.34	3 (0%) 85 83	45, 74, 120, 155	0
1	D	525/567 (92%)	-0.38	1 (0%) 91 90	44, 75, 121, 167	0
1	E	488/567 (86%)	0.18	7 (1%) 73 71	101, 118, 147, 190	0
1	F	490/567 (86%)	0.18	10 (2%) 65 62	102, 118, 151, 178	0
All	All	3082/3402 (90%)	-0.28	21 (0%) 84 82	41, 77, 136, 190	0

The worst 5 of 21 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	123	ILE	3.5
1	E	450	THR	3.3
1	E	363	LEU	2.8
1	E	421	GLY	2.7
1	E	132	ILE	2.5

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	GOL	F	601	6/6	0.95	0.07	74,86,91,92	0
2	GOL	B	602	6/6	0.97	0.09	71,75,80,82	0
2	GOL	B	601	6/6	0.98	0.10	51,61,64,69	0
2	GOL	A	601	6/6	0.98	0.08	45,51,55,58	0
2	GOL	C	601	6/6	0.98	0.09	59,62,66,75	0
2	GOL	C	602	6/6	0.98	0.08	50,52,63,68	0
2	GOL	D	601	6/6	0.98	0.08	51,57,61,68	0
2	GOL	A	602	6/6	0.98	0.09	71,79,86,91	0
2	GOL	D	602	6/6	0.99	0.06	40,52,58,60	0
2	GOL	A	603	6/6	0.99	0.07	59,70,74,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.