



# wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 22, 2020 – 12:52 PM BST

PDB ID : 5GKD  
Title : Structure of PL6 family alginate lyase AlyGC  
Authors : Zhang, Y.Z.; Wang, P.; Xu, F.  
Deposited on : 2016-07-04  
Resolution : 2.19 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.13.1

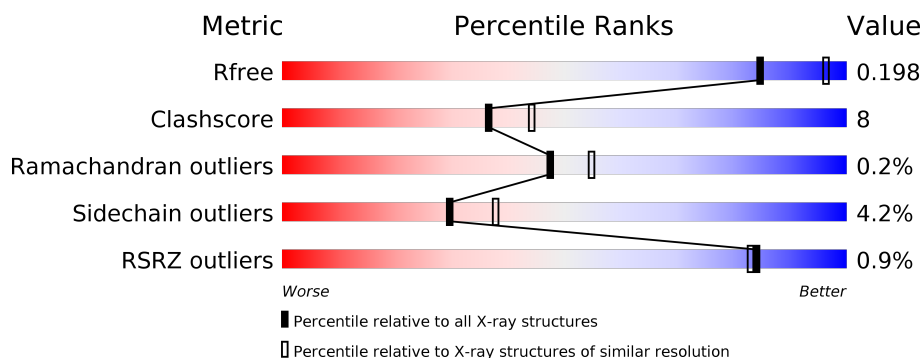
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.19 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))
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	726	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>84%</span> <span>14%</span> <span>•</span> </div> </div>
1	B	726	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>82%</span> <span>16%</span> <span>•</span> </div> </div>
1	C	726	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>83%</span> <span>15%</span> <span>•</span> </div> </div>
1	D	726	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> <span>%</span> <span>82%</span> <span>17%</span> <span>•</span> </div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	PO4	B	803	-	X	X	-
4	PO4	D	803	-	-	X	-
4	PO4	D	805	-	-	X	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 25339 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 AlyGC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	726	Total	C	N	O	S	0	0	0
			5667	3540	1012	1106	9			
1	B	726	Total	C	N	O	S	0	0	0
			5667	3540	1012	1106	9			
1	C	726	Total	C	N	O	S	0	0	0
			5667	3540	1012	1106	9			
1	D	726	Total	C	N	O	S	0	0	0
			5667	3540	1012	1106	9			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

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

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



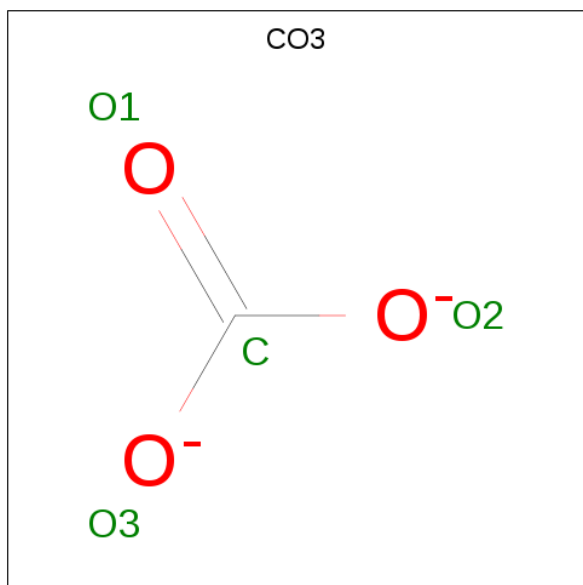
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	C	1	Total	O	P	0	0
			5	4	1		
4	C	1	Total	O	P	0	0
			5	4	1		
4	C	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	O	P	0	0
			5	4	1		
4	D	1	Total	O	P	0	0
			5	4	1		
4	D	1	Total	O	P	0	0
			5	4	1		

- Molecule 5 is CARBONATE ION (three-letter code: CO3) (formula: CO<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	1	3		
5	B	1	Total	C	O	0	0
			4	1	3		
5	C	1	Total	C	O	0	0
			4	1	3		
5	D	1	Total	C	O	0	0
			4	1	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	646	Total	O	0	0
			646	646		
6	B	693	Total	O	0	0
			693	693		

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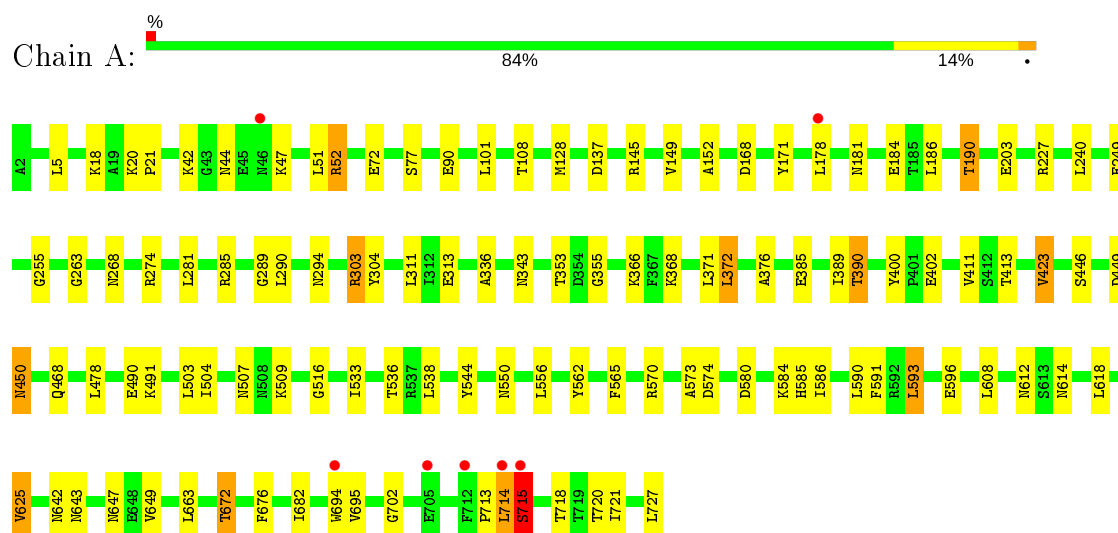
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	560	Total	O	0	0
			560	560		
6	D	596	Total	O	0	0
			596	596		



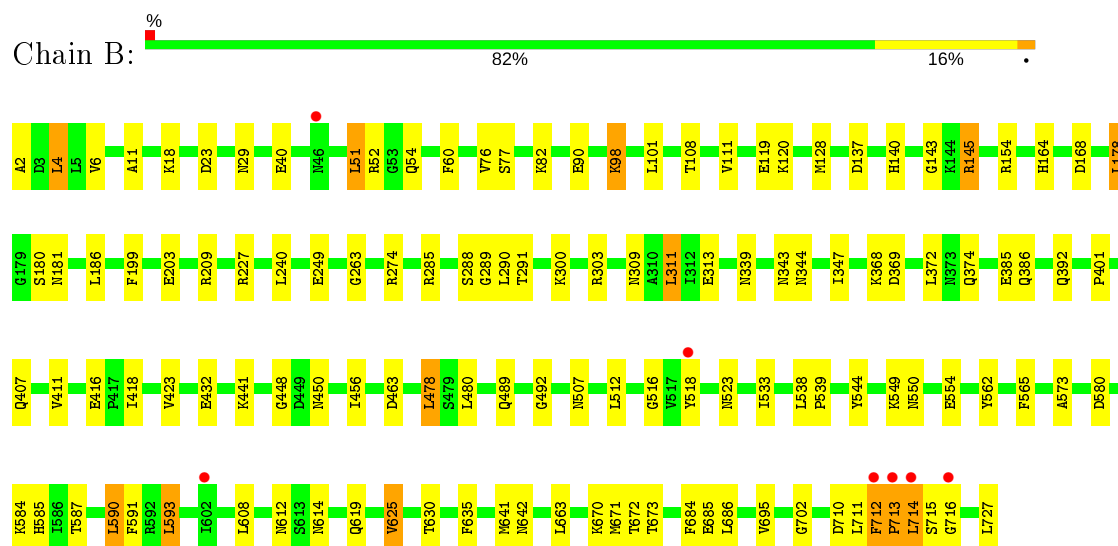
### 3 Residue-property plots

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.

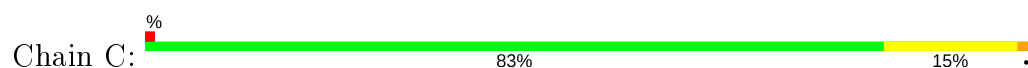
#### • Molecule 1: AlyGC

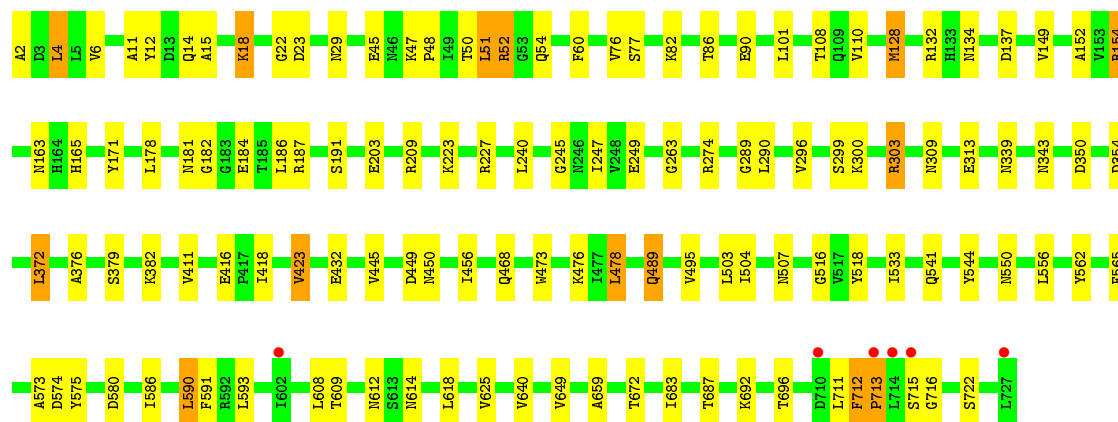


#### • Molecule 1: AlyGC

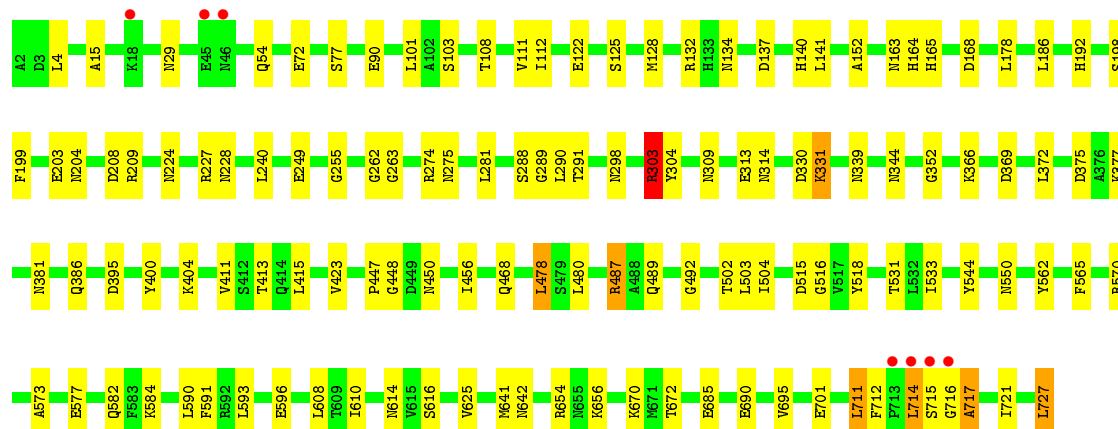
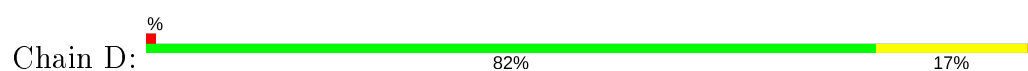


#### • Molecule 1: AlyGC





• Molecule 1: AlyGC



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	116.61Å 142.84Å 126.28Å 90.00° 111.09° 90.00°	Depositor
Resolution (Å)	34.45 – 2.19 49.91 – 2.19	Depositor EDS
% Data completeness (in resolution range)	98.4 (34.45-2.19) 98.4 (49.91-2.19)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.95 (at 2.20Å)	Xtriage
Refinement program	PHENIX (1.10.1 _2155: ???)	Depositor
R, $R_{free}$	0.165 , 0.197 0.166 , 0.198	Depositor DCC
$R_{free}$ test set	9768 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.7	Xtriage
Anisotropy	0.190	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 43.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	25339	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.21% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, CA, PO4, CO3

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	0.50	1/5776 (0.0%)	0.69	2/7830 (0.0%)
1	B	0.51	3/5776 (0.1%)	0.67	3/7830 (0.0%)
1	C	0.46	0/5776	0.64	0/7830
1	D	0.46	3/5776 (0.1%)	0.64	1/7830 (0.0%)
All	All	0.48	7/23104 (0.0%)	0.66	6/31320 (0.0%)

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	98	LYS	CE-NZ	-12.60	1.17	1.49
1	D	685	GLU	CD-OE1	-6.50	1.18	1.25
1	A	423	VAL	CB-CG1	-6.45	1.39	1.52
1	B	685	GLU	CD-OE2	-6.29	1.18	1.25
1	D	685	GLU	CD-OE2	-6.16	1.18	1.25

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	715	SER	CB-CA-C	-12.13	87.06	110.10
1	B	98	LYS	CD-CE-NZ	-11.91	84.31	111.70
1	A	715	SER	N-CA-C	10.94	140.54	111.00
1	B	686	LEU	CB-CA-C	6.59	122.72	110.20
1	D	303	ARG	N-CA-C	6.15	127.59	111.00

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	5667	0	5529	64	0
1	B	5667	0	5529	99	0
1	C	5667	0	5529	90	0
1	D	5667	0	5529	93	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	24	0	32	4	0
3	B	24	0	32	1	0
3	C	24	0	32	3	0
3	D	24	0	32	0	0
4	A	15	0	0	1	0
4	B	15	0	0	2	0
4	C	15	0	0	0	0
4	D	15	0	0	5	0
5	A	4	0	0	0	0
5	B	4	0	0	0	0
5	C	4	0	0	0	0
5	D	4	0	0	0	0
6	A	646	0	0	20	2
6	B	693	0	0	20	5
6	C	560	0	0	20	2
6	D	596	0	0	24	1
All	All	25339	0	22244	358	5

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 358 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:A:715:SER:HB2	6:A:913:HOH:O	1.11	1.24
1:A:446:SER:OG	1:A:468:GLN:OE1	1.74	1.04
1:C:712:PHE:O	6:C:902:HOH:O	1.78	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:350:ASP:OD2	6:C:903:HOH:O	1.84	0.95
1:B:714:LEU:H	1:B:714:LEU:HD23	1.31	0.94

All (5) 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)
6:B:1347:HOH:O	6:C:1189:HOH:O[2_647]	1.97	0.23
6:B:1235:HOH:O	6:D:1447:HOH:O[2_647]	2.01	0.19
6:A:1000:HOH:O	6:B:1561:HOH:O[2_556]	2.03	0.17
6:B:1434:HOH:O	6:C:1368:HOH:O[2_647]	2.17	0.03
6:A:1541:HOH:O	6:B:1590:HOH:O[2_556]	2.18	0.02

## 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	724/726 (100%)	702 (97%)	21 (3%)	1 (0%)	51	60
1	B	724/726 (100%)	707 (98%)	16 (2%)	1 (0%)	51	60
1	C	724/726 (100%)	698 (96%)	24 (3%)	2 (0%)	41	46
1	D	724/726 (100%)	707 (98%)	15 (2%)	2 (0%)	41	46
All	All	2896/2904 (100%)	2814 (97%)	76 (3%)	6 (0%)	47	55

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	713	PRO
1	C	715	SER
1	A	715	SER
1	D	717	ALA
1	D	448	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	621/621 (100%)	591 (95%)	30 (5%)	25	32
1	B	621/621 (100%)	594 (96%)	27 (4%)	29	36
1	C	621/621 (100%)	593 (96%)	28 (4%)	27	34
1	D	621/621 (100%)	602 (97%)	19 (3%)	40	51
All	All	2484/2484 (100%)	2380 (96%)	104 (4%)	30	38

5 of 104 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	489	GLN
1	C	18	LYS
1	D	489	GLN
1	B	507	ASN
1	B	608	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	351	GLN
1	C	468	GLN
1	D	165	HIS
1	C	165	HIS
1	C	614	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](#)

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 36 ligands modelled in this entry, 4 are monoatomic - leaving 32 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	$\# Z  > 2$
3	GOL	D	807	-	5,5,5	0.38	0	5,5,5	0.23	0
3	GOL	C	809	-	5,5,5	0.41	0	5,5,5	0.41	0
3	GOL	A	809	-	5,5,5	0.48	0	5,5,5	0.47	0
3	GOL	B	809	-	5,5,5	0.34	0	5,5,5	0.35	0
3	GOL	A	807	-	5,5,5	0.49	0	5,5,5	0.71	0
5	CO3	B	806	-	0,3,3	0.00	-	0,3,3	0.00	-
3	GOL	D	802	-	5,5,5	0.42	0	5,5,5	0.89	0
3	GOL	B	802	-	5,5,5	0.42	0	5,5,5	1.02	0
5	CO3	A	806	-	0,3,3	0.00	-	0,3,3	0.00	-
3	GOL	C	802	-	5,5,5	0.38	0	5,5,5	0.52	0
4	PO4	C	805	-	4,4,4	0.93	0	6,6,6	0.64	0
3	GOL	B	807	-	5,5,5	0.40	0	5,5,5	0.44	0
3	GOL	D	809	-	5,5,5	0.43	0	5,5,5	0.32	0
3	GOL	C	807	-	5,5,5	0.39	0	5,5,5	0.16	0
4	PO4	C	803	-	4,4,4	0.94	0	6,6,6	0.70	0
4	PO4	A	803	-	4,4,4	0.87	0	6,6,6	0.55	0
4	PO4	A	805	-	4,4,4	0.95	0	6,6,6	0.45	0
4	PO4	A	804	2	4,4,4	1.19	1 (25%)	6,6,6	1.36	1 (16%)
3	GOL	C	808	-	5,5,5	0.35	0	5,5,5	0.61	0
3	GOL	D	808	-	5,5,5	0.34	0	5,5,5	0.13	0
3	GOL	B	808	-	5,5,5	0.43	0	5,5,5	0.31	0
5	CO3	C	806	-	0,3,3	0.00	-	0,3,3	0.00	-
4	PO4	D	805	-	4,4,4	0.92	0	6,6,6	0.48	0
4	PO4	B	805	-	4,4,4	0.97	0	6,6,6	0.69	0
3	GOL	A	808	-	5,5,5	0.29	0	5,5,5	0.38	0
4	PO4	B	804	2	4,4,4	1.57	1 (25%)	6,6,6	1.08	0
4	PO4	D	803	-	4,4,4	0.78	0	6,6,6	0.60	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	CO3	D	806	-	0,3,3	0.00	-	0,3,3	0.00	-
4	PO4	C	804	2	4,4,4	1.11	0	6,6,6	1.29	1 (16%)
4	PO4	B	803	-	4,4,4	1.44	1 (25%)	6,6,6	3.05	5 (83%)
4	PO4	D	804	2	4,4,4	1.06	0	6,6,6	1.60	2 (33%)
3	GOL	A	802	-	5,5,5	0.39	0	5,5,5	0.88	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
3	GOL	D	808	-	-	4/4/4/4	-
3	GOL	B	808	-	-	2/4/4/4	-
3	GOL	B	807	-	-	4/4/4/4	-
3	GOL	D	807	-	-	2/4/4/4	-
3	GOL	C	809	-	-	4/4/4/4	-
3	GOL	D	809	-	-	0/4/4/4	-
3	GOL	A	809	-	-	3/4/4/4	-
3	GOL	B	809	-	-	4/4/4/4	-
3	GOL	A	807	-	-	2/4/4/4	-
3	GOL	C	807	-	-	2/4/4/4	-
3	GOL	C	808	-	-	2/4/4/4	-
3	GOL	D	802	-	-	4/4/4/4	-
3	GOL	A	808	-	-	2/4/4/4	-
3	GOL	B	802	-	-	4/4/4/4	-
3	GOL	A	802	-	-	2/4/4/4	-
3	GOL	C	802	-	-	2/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	803	PO4	P-O4	-2.44	1.47	1.54
4	A	804	PO4	P-O2	-2.12	1.48	1.54
4	B	804	PO4	P-O1	2.04	1.55	1.50

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	803	PO4	O4-P-O3	-5.38	90.70	107.97
4	B	803	PO4	O3-P-O2	3.34	118.70	107.97
4	A	804	PO4	O4-P-O3	2.54	116.11	107.97
4	D	804	PO4	O4-P-O1	2.37	119.56	110.89
4	D	804	PO4	O3-P-O1	-2.30	102.49	110.89

There are no chirality outliers.

5 of 43 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	807	GOL	O1-C1-C2-C3
3	C	809	GOL	O1-C1-C2-C3
3	A	809	GOL	O1-C1-C2-C3
3	B	809	GOL	C1-C2-C3-O3
3	D	802	GOL	O1-C1-C2-C3

There are no ring outliers.

9 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	809	GOL	1	0
3	B	809	GOL	1	0
3	A	807	GOL	2	0
3	C	802	GOL	2	0
4	A	803	PO4	1	0
4	D	805	PO4	2	0
3	A	808	GOL	2	0
4	D	803	PO4	3	0
4	B	803	PO4	2	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	726/726 (100%)	-0.33	7 (0%) 82 81	16, 24, 39, 89	0
1	B	726/726 (100%)	-0.33	7 (0%) 82 81	17, 24, 38, 74	0
1	C	726/726 (100%)	-0.21	6 (0%) 86 85	18, 28, 45, 84	0
1	D	726/726 (100%)	-0.21	7 (0%) 82 81	18, 28, 45, 90	0
All	All	2904/2904 (100%)	-0.27	27 (0%) 84 83	16, 26, 43, 90	0

The worst 5 of 27 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	714	LEU	11.3
1	A	714	LEU	6.6
1	B	714	LEU	6.3
1	D	715	SER	5.9
1	A	715	SER	4.9

### 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 monosaccharides 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, 95<sup>th</sup> 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	D	809	6/6	0.76	0.15	36,40,43,47	0
5	CO3	A	806	4/4	0.77	0.22	22,23,25,38	0
4	PO4	D	803	5/5	0.78	0.28	65,72,77,88	0
3	GOL	D	808	6/6	0.79	0.24	37,41,48,56	0
3	GOL	C	807	6/6	0.79	0.18	38,46,51,59	0
3	GOL	C	808	6/6	0.81	0.23	33,37,42,50	0
3	GOL	D	807	6/6	0.82	0.21	34,38,38,48	0
5	CO3	C	806	4/4	0.83	0.18	24,25,28,38	0
5	CO3	B	806	4/4	0.85	0.20	22,22,23,37	0
3	GOL	A	809	6/6	0.85	0.13	34,39,42,43	0
3	GOL	C	809	6/6	0.86	0.12	31,35,38,46	0
3	GOL	C	802	6/6	0.86	0.23	29,43,51,57	0
3	GOL	B	807	6/6	0.87	0.19	30,35,41,51	0
3	GOL	B	802	6/6	0.88	0.17	23,34,37,40	0
5	CO3	D	806	4/4	0.88	0.13	25,26,27,37	0
3	GOL	B	808	6/6	0.90	0.23	38,41,48,49	0
3	GOL	A	808	6/6	0.91	0.20	30,34,36,48	0
3	GOL	B	809	6/6	0.93	0.09	28,30,39,47	0
3	GOL	A	807	6/6	0.93	0.16	23,29,34,36	0
3	GOL	A	802	6/6	0.93	0.13	20,30,35,36	0
3	GOL	D	802	6/6	0.94	0.13	31,34,38,39	0
4	PO4	B	803	5/5	0.95	0.16	43,45,49,51	0
4	PO4	C	804	5/5	0.95	0.18	23,24,29,32	0
4	PO4	C	803	5/5	0.96	0.12	42,43,44,45	0
4	PO4	B	805	5/5	0.96	0.17	46,50,59,64	0
4	PO4	B	804	5/5	0.97	0.13	20,20,24,31	0
2	CA	C	801	1/1	0.97	0.16	22,22,22,22	0
4	PO4	D	805	5/5	0.97	0.22	47,48,51,65	0
4	PO4	A	803	5/5	0.97	0.09	33,36,45,47	0
4	PO4	A	804	5/5	0.97	0.15	17,20,23,27	0
4	PO4	C	805	5/5	0.97	0.15	40,51,55,60	0
4	PO4	D	804	5/5	0.98	0.12	24,25,27,31	0
2	CA	D	801	1/1	0.99	0.12	20,20,20,20	0
4	PO4	A	805	5/5	0.99	0.23	35,44,50,50	0
2	CA	B	801	1/1	1.00	0.14	19,19,19,19	0
2	CA	A	801	1/1	1.00	0.18	22,22,22,22	0

## 6.5 Other polymers

There are no such residues in this entry.