



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2025 – 09:26 PM EDT

PDB ID : 3MNM  
Title : Crystal structure of GAE domain of GGA2p from *Saccharomyces cerevisiae*  
Authors : Fang, P.; Wang, J.; Li, X.; Niu, L.; Teng, M.  
Deposited on : 2010-04-21  
Resolution : 1.73 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](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.

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 : 4.02b-467  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 1.21  
EDS : 3.0  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.004 (Gargrove)  
Density-Fitness : 1.0.11  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.41.4

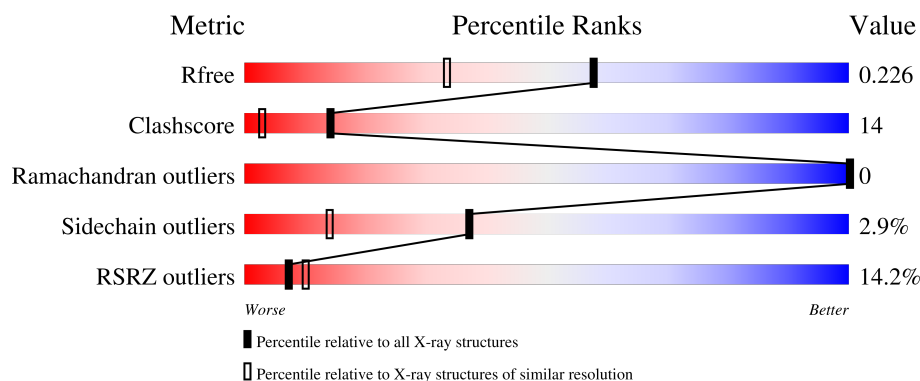
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.73 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))
$R_{free}$	164625	1043 (1.74-1.74)
Clashscore	180529	1119 (1.74-1.74)
Ramachandran outliers	177936	1112 (1.74-1.74)
Sidechain outliers	177891	1112 (1.74-1.74)
RSRZ outliers	164620	1043 (1.74-1.74)

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  $\geq 3$ , 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	123	<div> <div>13%</div> <div>72%</div> <div>17%</div> <div>9%</div> </div>
1	B	123	<div> <div>10%</div> <div>74%</div> <div>16%</div> <div>9%</div> </div>
1	C	123	<div> <div>15%</div> <div>75%</div> <div>13%</div> <div>11%</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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	586	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2965 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 ADP-ribosylation factor-binding protein GGA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	112	Total	C	N	O	S	0	6	0
			880	567	142	169	2			
1	B	112	Total	C	N	O	S	0	4	0
			877	564	142	169	2			
1	C	110	Total	C	N	O	S	0	1	0
			841	538	137	164	2			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	463	SER	-	expression tag	UNP P38817
A	464	LEU	-	expression tag	UNP P38817
B	463	SER	-	expression tag	UNP P38817
B	464	LEU	-	expression tag	UNP P38817
C	463	SER	-	expression tag	UNP P38817
C	464	LEU	-	expression tag	UNP P38817

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

*Continued on next page...*

*Continued from previous page...*

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

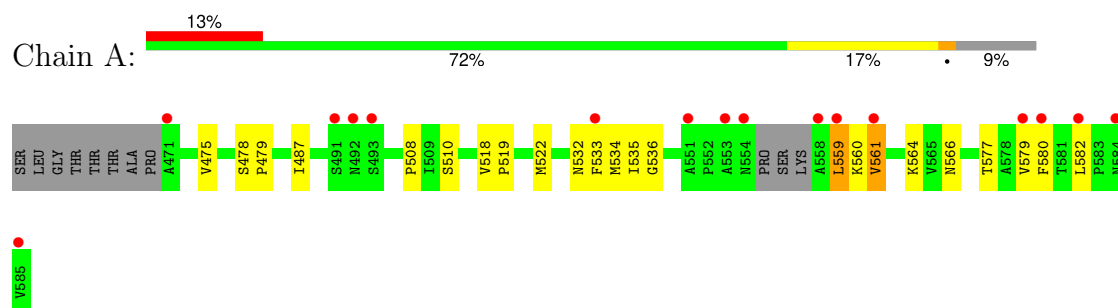
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	106	Total 106	O 106	0	0
4	B	110	Total 110	O 110	0	0
4	C	106	Total 106	O 106	0	0

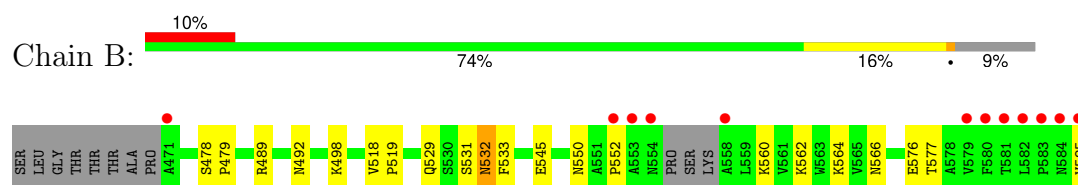
### 3 Residue-property plots [i](#)

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.

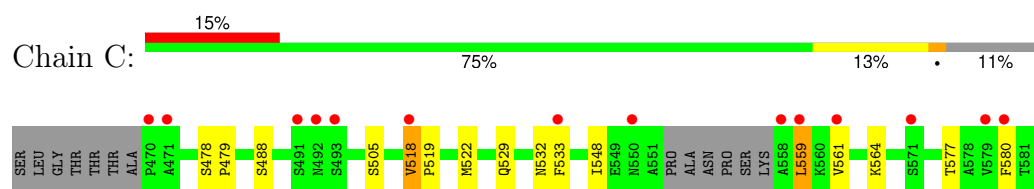
#### • Molecule 1: ADP-ribosylation factor-binding protein GGA2



#### • Molecule 1: ADP-ribosylation factor-binding protein GGA2



#### • Molecule 1: ADP-ribosylation factor-binding protein GGA2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	89.72Å 89.72Å 117.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.73 50.00 – 1.73	Depositor EDS
% Data completeness (in resolution range)	100.0 (50.00-1.73) 100.0 (50.00-1.73)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.18 (at 1.73Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, $R_{free}$	0.197 , 0.229 0.195 , 0.226	Depositor DCC
$R_{free}$ test set	2569 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	19.3	Xtriage
Anisotropy	0.002	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 52.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	2965	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.58% 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: MLY, SO4, 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	0.39	0/880	0.55	0/1195
1	B	0.39	0/869	0.56	0/1180
1	C	0.33	0/830	0.51	0/1128
All	All	0.38	0/2579	0.54	0/3503

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 [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	880	0	894	37	0
1	B	877	0	880	21	0
1	C	841	0	823	17	0
2	A	5	0	0	0	0
2	B	5	0	0	1	0
2	C	5	0	0	0	0
3	A	12	0	16	14	0
3	B	12	0	16	1	0
3	C	6	0	8	0	0
4	A	106	0	0	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	110	0	0	3	0
4	C	106	0	0	2	0
All	All	2965	0	2637	74	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 14.

All (74) 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:C:519:PRO:HD2	1:C:522:MET:HE2	1.40	0.99
1:A:508:PRO:HG3	3:A:586:GOL:H2	1.53	0.91
1:A:559:LEU:HG	1:A:582:LEU:HD22	1.55	0.89
1:C:522:MET:SD	1:C:559:LEU:HD21	2.13	0.88
1:A:535[B]:ILE:HG13	3:A:586:GOL:O2	1.77	0.85
1:C:522:MET:HE1	1:C:559:LEU:HD11	1.59	0.84
1:A:519:PRO:HD2	1:A:522:MET:HE2	1.59	0.83
1:A:508:PRO:HG3	3:A:586:GOL:C2	2.08	0.82
1:A:535[B]:ILE:CG1	3:A:586:GOL:H12	2.14	0.77
1:B:562:LYS:CG	3:B:3:GOL:H2	2.18	0.73
1:A:535[B]:ILE:HG13	3:A:586:GOL:H12	1.71	0.71
1:A:522:MET:HE1	1:A:559:LEU:HD11	1.73	0.70
1:A:519:PRO:HD2	1:A:522:MET:CE	2.23	0.67
1:B:560:MLY:HH13	2:B:1:SO4:O2	1.94	0.66
1:A:508:PRO:HB2	1:A:533[A]:PHE:CZ	2.31	0.66
1:B:564:LYS:HE2	1:B:566:ASN:OD1	1.98	0.64
1:A:508:PRO:CG	3:A:586:GOL:H2	2.25	0.64
1:A:510:SER:HB3	1:A:533[A]:PHE:CE1	2.33	0.64
1:A:535[B]:ILE:HG12	3:A:586:GOL:H12	1.80	0.63
1:B:498:MLY:HH22	4:B:589:HOH:O	1.98	0.62
1:C:519:PRO:HD2	1:C:522:MET:CE	2.25	0.60
1:A:533[A]:PHE:CD2	1:A:534:MET:N	2.70	0.60
1:B:550:ASN:HB3	4:B:317:HOH:O	2.03	0.59
1:C:488:SER:HA	1:C:585:VAL:HG11	1.83	0.59
1:B:564:LYS:HG3	1:B:577:THR:HG22	1.84	0.59
1:A:559:LEU:HB3	1:A:582:LEU:HB2	1.85	0.57
1:A:508:PRO:HG3	3:A:586:GOL:O2	2.06	0.56
1:A:508:PRO:CB	1:A:533[A]:PHE:CZ	2.89	0.55
1:B:498:MLY:HD3	1:B:545:GLU:HG2	1.89	0.55
1:A:564:LYS:HG3	1:A:577:THR:HG22	1.89	0.54
1:A:510:SER:HB3	1:A:533[A]:PHE:CD1	2.42	0.54

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:533[A]:PHE:CD2	4:C:364:HOH:O	2.54	0.53
1:A:518:VAL:HB	1:A:522:MET:CE	2.39	0.52
1:A:536:GLY:H	3:A:586:GOL:H12	1.74	0.52
1:A:561:VAL:HG12	1:A:580:PHE:HB3	1.92	0.52
1:A:535[B]:ILE:HG13	3:A:586:GOL:C1	2.40	0.52
1:A:508:PRO:HB3	1:A:533[A]:PHE:CE2	2.46	0.51
1:B:529:GLN:NE2	4:B:154:HOH:O	2.42	0.51
1:C:518:VAL:HG13	1:C:522:MET:HE2	1.92	0.51
1:A:508:PRO:HD3	3:A:586:GOL:H2	1.93	0.51
1:A:508:PRO:CD	3:A:586:GOL:H2	2.41	0.51
1:B:478:SER:HB3	1:B:576[B]:GLU:OE2	2.11	0.50
1:B:489:ARG:HB2	1:B:585:VAL:HG12	1.94	0.49
1:B:478:SER:HB2	1:B:479:PRO:HD2	1.94	0.49
1:B:478:SER:HB2	1:B:479:PRO:CD	2.42	0.49
1:A:519:PRO:HD3	1:A:560:MLY:O	2.11	0.49
1:B:532:ASN:HD22	1:B:533[B]:PHE:HD2	1.60	0.49
1:C:478:SER:HB2	1:C:479:PRO:HD2	1.94	0.49
1:A:522:MET:HE1	1:A:559:LEU:CD1	2.40	0.49
1:C:522:MET:SD	1:C:559:LEU:CD2	2.97	0.48
1:B:492:ASN:O	1:B:552:PRO:HA	2.13	0.48
1:B:518:VAL:HB	1:B:519:PRO:HD2	1.96	0.47
1:A:564:LYS:HE2	1:A:566:ASN:OD1	2.15	0.47
1:C:478:SER:HB2	1:C:479:PRO:CD	2.45	0.46
1:C:522:MET:HE3	1:C:548:ILE:HG21	1.97	0.46
1:A:536:GLY:H	3:A:586:GOL:C1	2.29	0.45
1:A:478:SER:HB2	1:A:479:PRO:CD	2.46	0.45
1:C:522:MET:HE3	1:C:548:ILE:CG2	2.47	0.45
1:B:518:VAL:HB	1:B:519:PRO:CD	2.47	0.44
1:A:533[A]:PHE:HD2	1:A:534:MET:N	2.15	0.44
1:C:518:VAL:HG13	1:C:522:MET:HB2	2.00	0.44
1:A:478:SER:HB2	1:A:479:PRO:HD2	2.00	0.44
1:B:532:ASN:HD22	1:B:532:ASN:H	1.66	0.44
1:A:487:ILE:HD13	1:A:582:LEU:HG	1.98	0.44
1:C:529:GLN:NE2	4:C:152:HOH:O	2.52	0.43
1:C:559:LEU:HG	1:C:582:LEU:HD12	2.00	0.42
1:C:561:VAL:HG22	1:C:580:PHE:O	2.18	0.42
1:B:531:SER:OG	1:B:533[B]:PHE:CE2	2.72	0.42
1:A:535[B]:ILE:HG13	3:A:586:GOL:C2	2.50	0.41
1:A:561:VAL:O	1:A:579:VAL:HA	2.20	0.41
1:C:564:LYS:HG3	1:C:577:THR:HG22	2.02	0.40
1:A:475:VAL:O	1:B:479:PRO:HD3	2.22	0.40

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	111/123 (90%)	110 (99%)	1 (1%)	0	100	100
1	B	109/123 (89%)	105 (96%)	4 (4%)	0	100	100
1	C	104/123 (85%)	103 (99%)	1 (1%)	0	100	100
All	All	324/369 (88%)	318 (98%)	6 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	98/107 (92%)	95 (97%)	3 (3%)	35	12
1	B	96/107 (90%)	95 (99%)	1 (1%)	73	61
1	C	91/107 (85%)	87 (96%)	4 (4%)	24	5
All	All	285/321 (89%)	277 (97%)	8 (3%)	37	15

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

Mol	Chain	Res	Type
1	A	532	ASN
1	A	559	LEU

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	561	VAL
1	B	532	ASN
1	C	505	SER
1	C	518	VAL
1	C	532	ASN
1	C	559	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	529	GLN
1	A	532	ASN
1	B	529	GLN
1	B	532	ASN
1	C	529	GLN
1	C	532	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

9 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	C	525	1	7,8,11	0.55	0	3,8,13	0.33	0
1	MLY	C	498	1	9,10,11	0.58	0	6,11,13	0.67	0
1	MLY	A	498	1	9,10,11	0.59	0	6,11,13	0.78	0
1	MLY	A	525	1	9,10,11	0.79	0	6,11,13	0.82	0
1	MLY	A	560	1	7,8,11	0.48	0	3,8,13	0.67	0
1	MLY	B	498	1	9,10,11	0.61	0	6,11,13	0.58	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	MLY	B	525	1	7,8,11	0.53	0	3,8,13	0.44	0
1	MLY	C	560	1	5,6,11	0.54	0	2,6,13	0.10	0
1	MLY	B	560	1	9,10,11	0.58	0	6,11,13	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	C	525	1	-	1/6/7/11	-
1	MLY	C	498	1	-	0/8/9/11	-
1	MLY	A	498	1	-	0/8/9/11	-
1	MLY	A	525	1	-	1/8/9/11	-
1	MLY	A	560	1	-	0/6/7/11	-
1	MLY	B	498	1	-	2/8/9/11	-
1	MLY	B	525	1	-	0/6/7/11	-
1	MLY	C	560	1	-	1/4/5/11	-
1	MLY	B	560	1	-	2/8/9/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	525	MLY	O-C-CA-CB
1	C	525	MLY	O-C-CA-CB
1	B	498	MLY	CD-CE-NZ-CH2
1	B	560	MLY	CG-CD-CE-NZ
1	B	560	MLY	CE-CD-CG-CB
1	B	498	MLY	CG-CD-CE-NZ
1	C	560	MLY	CA-CB-CG-CD

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	560	MLY	1	0
1	B	498	MLY	2	0

*Continued on next page...*

Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	560	MLY	1	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

8 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	$\# Z  > 2$
3	GOL	C	2	-	5,5,5	0.40	0	5,5,5	0.34	0
3	GOL	A	1	-	5,5,5	0.45	0	5,5,5	0.69	0
3	GOL	B	4	-	5,5,5	0.39	0	5,5,5	0.44	0
3	GOL	B	3	-	5,5,5	0.38	0	5,5,5	0.37	0
3	GOL	A	586	-	5,5,5	0.61	0	5,5,5	0.64	0
2	SO4	C	3	-	4,4,4	0.24	0	6,6,6	0.12	0
2	SO4	A	2	-	4,4,4	0.24	0	6,6,6	0.13	0
2	SO4	B	1	-	4,4,4	0.23	0	6,6,6	0.10	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	C	2	-	-	2/4/4/4	-
3	GOL	A	1	-	-	3/4/4/4	-
3	GOL	B	4	-	-	0/4/4/4	-
3	GOL	B	3	-	-	2/4/4/4	-
3	GOL	A	586	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	586	GOL	O1-C1-C2-O2
3	A	586	GOL	O1-C1-C2-C3
3	B	3	GOL	O1-C1-C2-C3
3	C	2	GOL	O1-C1-C2-C3
3	A	1	GOL	O1-C1-C2-C3
3	A	1	GOL	C1-C2-C3-O3
3	B	3	GOL	O1-C1-C2-O2
3	C	2	GOL	O1-C1-C2-O2
3	A	1	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	3	GOL	1	0
3	A	586	GOL	14	0
2	B	1	SO4	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

### 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, 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	109/123 (88%)	0.34	16 (14%) <b>7</b> <b>9</b>	8, 17, 47, 61	7 (6%)
1	B	109/123 (88%)	0.33	12 (11%) <b>12</b> <b>17</b>	7, 19, 45, 57	7 (6%)
1	C	107/123 (86%)	0.53	18 (16%) <b>5</b> <b>6</b>	12, 24, 50, 64	1 (0%)
All	All	325/369 (88%)	0.40	46 (14%) <b>7</b> <b>10</b>	7, 21, 47, 64	15 (4%)

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	470	PRO	5.8
1	A	580	PHE	5.1
1	B	580	PHE	5.1
1	A	579	VAL	4.8
1	B	471	ALA	4.4
1	C	585	VAL	4.3
1	B	554	ASN	4.0
1	B	581	THR	3.9
1	C	579	VAL	3.7
1	B	585	VAL	3.7
1	C	584	ASN	3.6
1	A	533[A]	PHE	3.4
1	A	559	LEU	3.4
1	A	561	VAL	3.3
1	B	553	ALA	3.3
1	C	559	LEU	3.3
1	A	558	ALA	3.2
1	A	554	ASN	3.1
1	C	492	ASN	3.1
1	C	582	LEU	3.1
1	C	583	PRO	3.1
1	C	580	PHE	3.1
1	A	553	ALA	3.0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	584	ASN	2.8
1	C	493	SER	2.7
1	A	471	ALA	2.7
1	C	533[A]	PHE	2.6
1	A	491	SER	2.6
1	C	550	ASN	2.5
1	C	491	SER	2.4
1	B	552	PRO	2.4
1	A	585	VAL	2.4
1	B	582	LEU	2.4
1	A	492	ASN	2.3
1	A	493	SER	2.3
1	A	582	LEU	2.2
1	B	584	ASN	2.2
1	A	551	ALA	2.1
1	C	558	ALA	2.1
1	B	583	PRO	2.1
1	C	561	VAL	2.1
1	B	558	ALA	2.1
1	B	579	VAL	2.1
1	C	471	ALA	2.1
1	C	571	SER	2.1
1	C	518	VAL	2.1

## 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, 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(Å <sup>2</sup> )	Q<0.9
1	MLY	C	560	7/12	0.82	0.17	27,32,42,46	0
1	MLY	A	560	9/12	0.89	0.12	24,26,37,45	0
1	MLY	B	525	9/12	0.92	0.12	16,19,45,53	0
1	MLY	B	560	11/12	0.94	0.12	20,26,40,50	0
1	MLY	C	525	9/12	0.94	0.09	13,15,32,33	0
1	MLY	B	498	11/12	0.94	0.09	15,17,33,39	0
1	MLY	A	498	11/12	0.95	0.09	14,16,21,27	0
1	MLY	A	525	11/12	0.95	0.11	13,15,37,37	0
1	MLY	C	498	11/12	0.97	0.07	16,20,27,33	0

### 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
2	SO4	C	3	5/5	0.80	0.12	54,64,73,78	0
3	GOL	A	586	6/6	0.82	0.35	19,29,31,37	0
3	GOL	B	3	6/6	0.86	0.12	27,39,42,51	0
2	SO4	A	2	5/5	0.89	0.10	44,45,51,55	0
3	GOL	B	4	6/6	0.90	0.11	18,28,29,36	0
3	GOL	A	1	6/6	0.92	0.11	14,23,26,27	0
3	GOL	C	2	6/6	0.93	0.10	20,28,31,36	0
2	SO4	B	1	5/5	0.95	0.10	33,40,50,51	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.