



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:36 AM GMT

PDB ID : 2XON  
Title : STRUCTURE OF TMCBM61 IN COMPLEX WITH BETA-1,4-GALACTO  
TRIOSE AT 1.4 Å RESOLUTION  
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Deposited on : 2010-08-20  
Resolution : 1.40 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  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

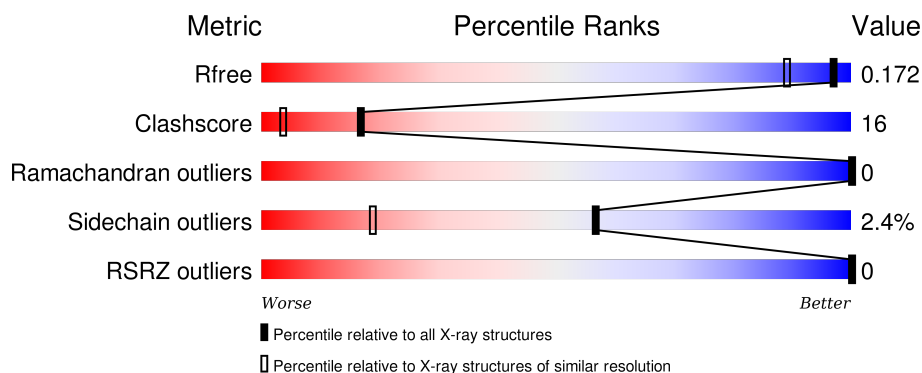
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.40 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))
$R_{free}$	91344	1199 (1.40-1.40)
Clashscore	102246	1295 (1.40-1.40)
Ramachandran outliers	100387	1259 (1.40-1.40)
Sidechain outliers	100360	1258 (1.40-1.40)
RSRZ outliers	91569	1198 (1.40-1.40)

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. 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	152	 82% 13% • 5%
1	L	152	 79% 16% • •

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	SO4	L	1171	-	X	-	-
4	EDO	A	1168	-	-	X	-
4	EDO	A	1169	-	-	X	-
4	EDO	A	1171	-	-	-	X
4	EDO	A	1172	-	-	-	X
4	EDO	A	1173	-	-	X	X
4	EDO	A	1183	-	-	X	-
4	EDO	L	1172	-	-	-	X
4	EDO	L	1173	-	-	-	X
4	EDO	L	1175	-	-	X	-
4	EDO	L	1176	-	-	X	-
4	EDO	L	1178	-	-	X	-
4	EDO	L	1179	-	-	X	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3225 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 ARABINOGALACTAN ENDO-1,4-BETA-GALACTOSIDASE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	145	Total	C	N	O	0	16	0
			1253	814	209	230			
1	L	146	Total	C	N	O	0	11	0
			1240	799	207	234			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	15	GLY	-	EXPRESSION TAG	UNP Q9X0S8
A	16	SER	-	EXPRESSION TAG	UNP Q9X0S8
A	17	HIS	-	EXPRESSION TAG	UNP Q9X0S8
A	18	MET	-	EXPRESSION TAG	UNP Q9X0S8
A	19	ALA	-	EXPRESSION TAG	UNP Q9X0S8
A	20	SER	-	EXPRESSION TAG	UNP Q9X0S8
L	15	GLY	-	EXPRESSION TAG	UNP Q9X0S8
L	16	SER	-	EXPRESSION TAG	UNP Q9X0S8
L	17	HIS	-	EXPRESSION TAG	UNP Q9X0S8
L	18	MET	-	EXPRESSION TAG	UNP Q9X0S8
L	19	ALA	-	EXPRESSION TAG	UNP Q9X0S8
L	20	SER	-	EXPRESSION TAG	UNP Q9X0S8

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

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

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	L	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	L	1	Total 4	C 2	O 2	0	0
4	L	1	Total 4	C 2	O 2	0	0
4	L	1	Total 4	C 2	O 2	0	0
4	L	1	Total 4	C 2	O 2	0	0
4	L	1	Total 4	C 2	O 2	0	0
4	L	1	Total 4	C 2	O 2	0	0
4	L	1	Total 3	C 2	O 1	0	0
4	L	1	Total 4	C 2	O 2	0	0

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

- Molecule 5 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	3	Total	C	O	0	0
			34	18	16		
5	L	3	Total	C	O	0	0
			34	18	16		


- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	285	Total	O	0	0
			285	285		
6	L	268	Total	O	0	0
			268	268		

### 3 Residue-property plots


These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

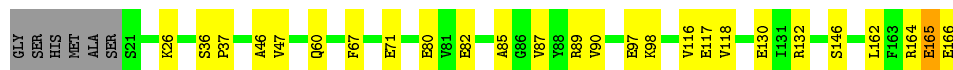
- Molecule 1: ARABINOGLACTAN ENDO-1,4-BETA-GALACTOSIDASE

Chain A: 



- Molecule 1: ARABINOGLACTAN ENDO-1,4-BETA-GALACTOSIDASE

Chain L: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	48.90Å 61.85Å 55.53Å 90.00° 105.14° 90.00°	Depositor
Resolution (Å)	34.26 – 1.40 34.26 – 1.40	Depositor EDS
% Data completeness (in resolution range)	95.0 (34.26-1.40) 95.1 (34.26-1.40)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.35 (at 1.40Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, $R_{free}$	0.116 , 0.167 0.130 , 0.172	Depositor DCC
$R_{free}$ test set	3009 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	12.5	Xtriage
Anisotropy	0.223	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 41.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 59810 reflections	Xtriage
$F_o, F_c$ correlation	0.98	EDS
Total number of atoms	3225	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.93% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, GAL, SO4, EDO

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	1.12	3/1333 (0.2%)	1.07	3/1802 (0.2%)
1	L	1.21	5/1301 (0.4%)	1.05	1/1756 (0.1%)
All	All	1.17	8/2634 (0.3%)	1.06	4/3558 (0.1%)

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	L	0	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	82	GLU	CD-OE2	-14.52	1.09	1.25
1	A	21	SER	CB-OG	6.87	1.51	1.42
1	A	114	ARG	CG-CD	-5.90	1.37	1.51
1	A	97	GLU	CD-OE1	-5.31	1.19	1.25
1	L	47	VAL	CB-CG2	-5.30	1.41	1.52
1	L	146	SER	CB-OG	-5.24	1.35	1.42
1	L	80[A]	GLU	CD-OE2	5.16	1.31	1.25
1	L	80[B]	GLU	CD-OE2	5.16	1.31	1.25

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	97	GLU	CA-C-N	-8.46	98.58	117.20
1	A	83[A]	LEU	CB-CA-C	-5.40	99.94	110.20
1	A	83[B]	LEU	CB-CA-C	-5.40	99.94	110.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	164	ARG	NE-CZ-NH1	5.39	123.00	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L	165	GLU	Peptide

## 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	1253	0	1262	20	2
1	L	1240	0	1213	40	0
2	A	1	0	0	0	0
2	L	1	0	0	0	0
3	A	5	0	0	0	0
3	L	5	0	0	0	0
4	A	60	0	90	26	0
4	L	39	0	57	30	2
5	A	34	0	30	2	0
5	L	34	0	30	2	0
6	A	285	0	0	20	0
6	L	268	0	0	16	0
All	All	3225	0	2682	88	2

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 16.

All (88) 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:L:90[A]:VAL:HG22	1:L:162:LEU:HD13	1.23	1.15
1:L:90[A]:VAL:CG2	1:L:162:LEU:HD13	1.85	1.06
1:A:21:SER:N	6:A:2004:HOH:O	1.87	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71[A]:GLU:OE1	6:A:2128:HOH:O	1.87	0.91
1:L:90[A]:VAL:HG22	1:L:162:LEU:CD1	2.00	0.91
1:L:166[B]:GLU:CA	1:L:166[B]:GLU:OE2	2.26	0.82
4:A:1168:EDO:H12	4:A:1169:EDO:C1	2.11	0.79
4:A:1175:EDO:H21	6:A:2267:HOH:O	1.84	0.77
4:L:1177:EDO:H21	6:L:2081:HOH:O	1.83	0.77
1:L:164:ARG:HE	4:L:1175:EDO:C2	1.98	0.77
1:A:21:SER:CA	6:A:2004:HOH:O	2.28	0.77
5:L:1169:GAL:O4	4:L:1178:EDO:H22	1.87	0.75
1:A:21:SER:HA	6:A:2004:HOH:O	1.86	0.74
1:L:164:ARG:HE	4:L:1175:EDO:H22	1.52	0.74
4:L:1178:EDO:C1	6:L:2251:HOH:O	2.36	0.74
1:A:90[A]:VAL:HG22	1:A:162:LEU:HD13	1.70	0.73
1:A:116[B]:VAL:HG21	6:A:2206:HOH:O	1.86	0.73
1:L:117[B]:GLU:OE1	6:L:2181:HOH:O	2.07	0.73
4:A:1183:EDO:H21	6:A:2275:HOH:O	1.88	0.71
1:L:87:VAL:HB	4:L:1175:EDO:H11	1.73	0.71
1:L:71:GLU:HG2	6:L:2134:HOH:O	1.90	0.70
4:L:1178:EDO:H21	6:L:2261:HOH:O	1.92	0.70
4:A:1183:EDO:H12	6:A:2282:HOH:O	1.90	0.70
1:L:166[A]:GLU:OE2	6:L:2240:HOH:O	2.09	0.69
4:A:1168:EDO:C1	4:A:1169:EDO:C1	2.71	0.69
4:A:1168:EDO:H12	4:A:1169:EDO:H11	1.73	0.69
1:A:116[B]:VAL:CG2	6:A:2206:HOH:O	2.41	0.68
4:A:1168:EDO:H12	4:A:1169:EDO:H12	1.75	0.68
4:L:1178:EDO:H12	6:L:2251:HOH:O	1.93	0.68
4:A:1183:EDO:C2	6:A:2275:HOH:O	2.42	0.68
1:L:132[B]:ARG:NH1	4:L:1179:EDO:H22	2.10	0.66
1:A:118[B]:VAL:HG22	6:A:2181:HOH:O	1.96	0.66
4:A:1176:EDO:O2	6:A:2270:HOH:O	2.15	0.65
4:A:1168:EDO:C1	4:A:1169:EDO:H12	2.27	0.65
1:L:118[B]:VAL:HG12	6:L:2186:HOH:O	1.96	0.64
1:A:39[B]:ARG:NE	4:A:1173:EDO:O1	2.31	0.64
4:L:1178:EDO:H11	6:L:2251:HOH:O	1.96	0.63
1:L:164:ARG:HH21	4:L:1175:EDO:H21	1.63	0.63
1:A:39[B]:ARG:HE	4:A:1173:EDO:C1	2.11	0.62
1:L:132[B]:ARG:HH12	4:L:1179:EDO:H22	1.65	0.61
1:L:164:ARG:NE	4:L:1175:EDO:H22	2.16	0.60
1:A:39[B]:ARG:NH2	4:A:1173:EDO:O1	2.35	0.58
4:A:1184:EDO:H12	6:A:2276:HOH:O	2.02	0.58
1:L:89:ARG:NH1	1:L:132[A]:ARG:HD3	2.19	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:87:VAL:H	4:L:1175:EDO:H22	1.67	0.58
1:L:90[A]:VAL:CG2	1:L:162:LEU:CD1	2.68	0.57
1:L:132[B]:ARG:HH12	4:L:1179:EDO:C2	2.19	0.56
1:A:39[B]:ARG:CZ	4:A:1173:EDO:O1	2.55	0.55
1:A:90[A]:VAL:CG2	1:A:162:LEU:HD13	2.36	0.54
1:L:89:ARG:NH2	4:L:1176:EDO:O2	2.40	0.54
1:A:106[A]:VAL:HG12	1:A:143:ILE:CD1	2.37	0.54
4:L:1173:EDO:H22	6:L:2157:HOH:O	2.07	0.53
1:L:36:SER:OG	4:L:1174:EDO:H21	2.09	0.53
1:L:165:GLU:O	4:L:1175:EDO:O1	2.25	0.53
1:L:89:ARG:HH21	4:L:1176:EDO:C2	2.22	0.52
1:A:54:PRO:HG3	4:A:1171:EDO:H12	1.92	0.51
1:A:117[B]:GLU:OE2	6:A:2182:HOH:O	2.20	0.50
1:L:89:ARG:HD2	4:L:1176:EDO:H21	1.93	0.49
1:L:37:PRO:HA	4:L:1174:EDO:H21	1.92	0.49
1:L:130:GLU:OE2	4:L:1176:EDO:H22	2.12	0.49
1:L:37:PRO:HA	4:L:1174:EDO:C2	2.43	0.48
5:A:1182:GAL:O4	4:A:1183:EDO:H11	2.13	0.48
1:L:164:ARG:HH21	4:L:1175:EDO:C2	2.27	0.48
1:L:71:GLU:CG	6:L:2134:HOH:O	2.56	0.48
5:L:1169:GAL:H2	4:L:1178:EDO:H11	1.95	0.47
1:L:85:ALA:HB2	6:L:2218:HOH:O	2.14	0.47
1:A:39[B]:ARG:HE	4:A:1173:EDO:H11	1.78	0.46
1:A:130:GLU:OE2	4:A:1176:EDO:C1	2.63	0.46
1:L:116:VAL:HG21	4:L:1179:EDO:H12	1.98	0.46
4:A:1183:EDO:C1	6:A:2282:HOH:O	2.55	0.46
4:A:1173:EDO:H12	6:A:2038:HOH:O	2.16	0.45
1:L:26:LYS:HE2	6:L:2014:HOH:O	2.17	0.45
1:L:117[B]:GLU:CD	6:L:2181:HOH:O	2.53	0.45
1:A:106[A]:VAL:HG12	1:A:143:ILE:HD12	2.00	0.44
1:L:130:GLU:OE1	1:L:132[A]:ARG:NE	2.38	0.43
4:A:1168:EDO:H11	4:A:1169:EDO:C1	2.49	0.43
4:A:1175:EDO:C2	6:A:2267:HOH:O	2.55	0.43
4:A:1184:EDO:H22	6:A:2284:HOH:O	2.19	0.43
5:A:1182:GAL:H2	4:A:1183:EDO:H22	2.01	0.42
4:L:1178:EDO:C2	6:L:2261:HOH:O	2.59	0.42
1:L:87:VAL:H	4:L:1175:EDO:C2	2.32	0.42
1:L:89:ARG:HH21	4:L:1176:EDO:H21	1.83	0.41
4:A:1185:EDO:O2	6:A:2276:HOH:O	2.15	0.41
1:L:46:ALA:O	1:L:67:PHE:HA	2.20	0.41
1:L:166[A]:GLU:HG3	6:L:2242:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117[B]:GLU:CD	6:A:2182:HOH:O	2.59	0.40

All (2) 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)
1:A:80:GLU:OE1	4:L:1180:EDO:O2[1_556]	1.88	0.32
1:A:142:LYS:NZ	4:L:1172:EDO:C1[2_656]	2.16	0.04

## 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	159/152 (105%)	158 (99%)	1 (1%)	0	100	100
1	L	154/152 (101%)	153 (99%)	1 (1%)	0	100	100
All	All	313/304 (103%)	311 (99%)	2 (1%)	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	139/128 (109%)	133 (96%)	6 (4%)	35	5

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	135/128 (106%)	132 (98%)	3 (2%)	60	22
All	All	274/256 (107%)	265 (97%)	9 (3%)	57	10

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45[A]	LYS
1	A	45[B]	LYS
1	A	60	GLN
1	A	98[A]	LYS
1	A	98[B]	LYS
1	A	164	ARG
1	L	60	GLN
1	L	98[A]	LYS
1	L	98[B]	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	ASN
1	A	128	ASN
1	L	66	ASN
1	L	128	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

6 carbohydrates 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	$\# Z  > 2$
5	GAL	A	1180	5	12,12,12	0.66	0	17,17,17	1.20	2 (11%)
5	GAL	A	1181	5	11,11,12	0.91	0	14,15,17	0.93	1 (7%)
5	GAL	A	1182	5	11,11,12	0.71	0	14,15,17	1.05	1 (7%)
5	GAL	L	1167	5	12,12,12	0.95	1 (8%)	17,17,17	1.14	1 (5%)
5	GAL	L	1168	5	11,11,12	0.61	0	14,15,17	1.08	1 (7%)
5	GAL	L	1169	5	11,11,12	0.66	0	14,15,17	1.22	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
5	GAL	A	1180	5	-	0/2/22/22	0/1/1/1
5	GAL	A	1181	5	-	0/2/19/22	0/1/1/1
5	GAL	A	1182	5	-	0/2/19/22	0/1/1/1
5	GAL	L	1167	5	-	0/2/22/22	0/1/1/1
5	GAL	L	1168	5	-	0/2/19/22	0/1/1/1
5	GAL	L	1169	5	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	L	1167	GAL	O5-C5	-2.30	1.38	1.44

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1180	GAL	C1-C2-C3	-2.86	106.17	110.43
5	A	1182	GAL	O2-C2-C1	-2.47	104.25	109.21
5	A	1181	GAL	C1-O5-C5	2.01	114.80	112.25
5	A	1180	GAL	O1-C1-C2	2.08	114.79	109.21
5	L	1168	GAL	O5-C5-C6	2.09	111.86	107.35
5	L	1167	GAL	O2-C2-C3	2.53	116.04	110.34

There are no chirality outliers.



There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1182	GAL	2	0
5	L	1169	GAL	2	0

## 5.6 Ligand geometry [i](#)

Of 29 ligands modelled in this entry, 2 are monoatomic - leaving 27 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	SO4	A	1167	-	4,4,4	1.89	2 (50%)	6,6,6	0.94	0
4	EDO	A	1168	-	3,3,3	0.61	0	2,2,2	0.84	0
4	EDO	A	1169	-	3,3,3	0.46	0	2,2,2	0.81	0
4	EDO	A	1170	-	3,3,3	0.38	0	2,2,2	0.26	0
4	EDO	A	1171	-	3,3,3	0.67	0	2,2,2	2.06	1 (50%)
4	EDO	A	1172	-	3,3,3	1.34	0	2,2,2	1.25	0
4	EDO	A	1173	-	3,3,3	0.62	0	2,2,2	0.50	0
4	EDO	A	1174	-	3,3,3	0.64	0	2,2,2	0.13	0
4	EDO	A	1175	-	3,3,3	0.31	0	2,2,2	0.66	0
4	EDO	A	1176	-	3,3,3	0.35	0	2,2,2	0.44	0
4	EDO	A	1177	-	3,3,3	0.59	0	2,2,2	0.56	0
4	EDO	A	1178	-	3,3,3	0.29	0	2,2,2	1.48	1 (50%)
4	EDO	A	1179	-	3,3,3	0.20	0	2,2,2	1.90	0
4	EDO	A	1183	-	3,3,3	0.38	0	2,2,2	0.31	0
4	EDO	A	1184	-	3,3,3	0.72	0	2,2,2	0.17	0
4	EDO	A	1185	-	3,3,3	0.25	0	2,2,2	0.23	0
3	SO4	L	1171	-	4,4,4	1.77	2 (50%)	6,6,6	1.35	2 (33%)
4	EDO	L	1172	-	3,3,3	0.23	0	2,2,2	0.65	0
4	EDO	L	1173	-	3,3,3	0.39	0	2,2,2	0.27	0
4	EDO	L	1174	-	3,3,3	0.85	0	2,2,2	1.04	0
4	EDO	L	1175	-	3,3,3	0.36	0	2,2,2	0.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	EDO	L	1176	-	3,3,3	0.27	0	2,2,2	1.85	1 (50%)
4	EDO	L	1177	-	2,2,3	0.44	0	1,1,2	0.92	0
4	EDO	L	1178	-	3,3,3	0.36	0	2,2,2	0.51	0
4	EDO	L	1179	-	3,3,3	0.35	0	2,2,2	0.31	0
4	EDO	L	1180	-	3,3,3	0.32	0	2,2,2	0.88	0
4	EDO	L	1181	-	3,3,3	0.55	0	2,2,2	0.94	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	SO4	A	1167	-	-	0/0/0/0	0/0/0/0
4	EDO	A	1168	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1169	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1170	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1171	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1172	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1173	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1174	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1175	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1176	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1177	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1178	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1179	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1183	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1184	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1185	-	-	0/1/1/1	0/0/0/0
3	SO4	L	1171	-	-	0/0/0/0	0/0/0/0
4	EDO	L	1172	-	-	0/1/1/1	0/0/0/0
4	EDO	L	1173	-	-	0/1/1/1	0/0/0/0
4	EDO	L	1174	-	-	0/1/1/1	0/0/0/0
4	EDO	L	1175	-	-	0/1/1/1	0/0/0/0
4	EDO	L	1176	-	-	0/1/1/1	0/0/0/0
4	EDO	L	1177	-	-	0/0/0/1	0/0/0/0
4	EDO	L	1178	-	-	0/1/1/1	0/0/0/0
4	EDO	L	1179	-	-	0/1/1/1	0/0/0/0
4	EDO	L	1180	-	-	0/1/1/1	0/0/0/0
4	EDO	L	1181	-	-	0/1/1/1	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1167	SO4	O4-S	-2.41	1.38	1.47
3	L	1171	SO4	O3-S	-2.06	1.40	1.47
3	L	1171	SO4	O1-S	2.57	1.55	1.47
3	A	1167	SO4	O3-S	2.78	1.57	1.47

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	1176	EDO	O1-C1-C2	-2.51	94.53	112.54
4	A	1171	EDO	O1-C1-C2	-2.39	95.40	112.54
3	L	1171	SO4	O2-S-O1	-2.36	102.03	109.50
4	A	1178	EDO	O1-C1-C2	-2.07	97.68	112.54
3	L	1171	SO4	O4-S-O3	-2.01	100.79	108.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

18 monomers are involved in 58 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1168	EDO	6	0
4	A	1169	EDO	6	0
4	A	1171	EDO	1	0
4	A	1173	EDO	6	0
4	A	1175	EDO	2	0
4	A	1176	EDO	2	0
4	A	1183	EDO	6	0
4	A	1184	EDO	2	0
4	A	1185	EDO	1	0
4	L	1172	EDO	0	1
4	L	1173	EDO	1	0
4	L	1174	EDO	3	0
4	L	1175	EDO	9	0
4	L	1176	EDO	5	0
4	L	1177	EDO	1	0
4	L	1178	EDO	7	0
4	L	1179	EDO	4	0
4	L	1180	EDO	0	1

## 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	145/152 (95%)	-0.48	0 100 100	9, 12, 19, 32	4 (2%)
1	L	146/152 (96%)	-0.56	0 100 100	9, 12, 19, 24	6 (4%)
All	All	291/304 (95%)	-0.52	0 100 100	9, 12, 19, 32	10 (3%)

There are no RSRZ outliers to report.

### 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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
5	GAL	L	1168	11/12	0.98	0.06	0.24	13,14,18,25	0
5	GAL	A	1182	11/12	0.99	0.07	0.21	9,10,11,13	0
5	GAL	A	1181	11/12	0.98	0.07	-0.27	12,13,15,22	0
5	GAL	L	1169	11/12	0.98	0.05	-0.68	11,12,13,13	0
5	GAL	A	1180	12/12	0.95	0.08	-	16,20,24,30	0
5	GAL	L	1167	12/12	0.91	0.10	-	19,24,31,35	0

## 6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	EDO	L	1173	4/4	0.83	0.22	18.07	12,14,14,19	4
4	EDO	A	1171	4/4	0.89	0.24	17.02	8,18,18,19	4
4	EDO	A	1173	4/4	0.72	0.27	7.16	32,34,38,42	1
4	EDO	L	1179	4/4	0.86	0.26	5.94	31,36,37,39	0
4	EDO	A	1172	4/4	0.92	0.09	3.89	24,28,30,30	0
4	EDO	L	1172	4/4	0.86	0.14	2.28	22,23,24,25	4
4	EDO	L	1175	4/4	0.75	0.27	1.97	38,41,42,44	0
3	SO4	L	1171	5/5	0.97	0.10	1.58	10,11,14,14	5
4	EDO	L	1174	4/4	0.68	0.15	1.49	35,36,38,39	0
3	SO4	A	1167	5/5	0.97	0.11	0.98	9,12,13,15	5
4	EDO	A	1184	4/4	0.86	0.18	0.42	31,33,33,36	0
4	EDO	L	1181	4/4	0.94	0.16	0.29	12,16,19,23	4
2	CA	L	1170	1/1	1.00	0.03	-1.93	11,11,11,11	0
2	CA	A	1166	1/1	1.00	0.03	-1.93	9,9,9,9	0
4	EDO	L	1180	4/4	0.85	0.22	-	45,45,45,48	0
4	EDO	A	1177	4/4	0.61	0.28	-	43,44,44,44	0
4	EDO	L	1178	4/4	0.95	0.18	-	26,31,32,33	0
4	EDO	A	1176	4/4	0.88	0.27	-	16,23,23,29	4
4	EDO	A	1183	4/4	0.92	0.27	-	18,19,22,23	4
4	EDO	A	1174	4/4	0.49	0.29	-	58,58,58,59	0
4	EDO	A	1185	4/4	0.84	0.15	-	18,21,21,21	4
4	EDO	A	1168	4/4	0.67	0.31	-	24,32,36,43	1
4	EDO	A	1175	4/4	0.94	0.14	-	20,21,22,24	4
4	EDO	L	1177	3/4	0.68	0.23	-	30,30,36,36	0
4	EDO	A	1169	4/4	0.87	0.11	-	17,19,19,24	4
4	EDO	A	1170	4/4	0.88	0.14	-	19,20,22,26	4
4	EDO	A	1178	4/4	0.90	0.11	-	25,28,29,33	0
4	EDO	A	1179	4/4	0.94	0.21	-	19,28,30,31	2
4	EDO	L	1176	4/4	0.88	0.21	-	18,21,22,23	4

## 6.5 Other polymers ⓘ

There are no such residues in this entry.