



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:00 PM UTC

PDB ID : 4HTY / pdb_00004hty
Title : Crystal Structure of a metagenome-derived cellulase Cel5A
Authors : Zhuang, N.; Lee, K.H.
Deposited on : 2012-11-02
Resolution : 2.00 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

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

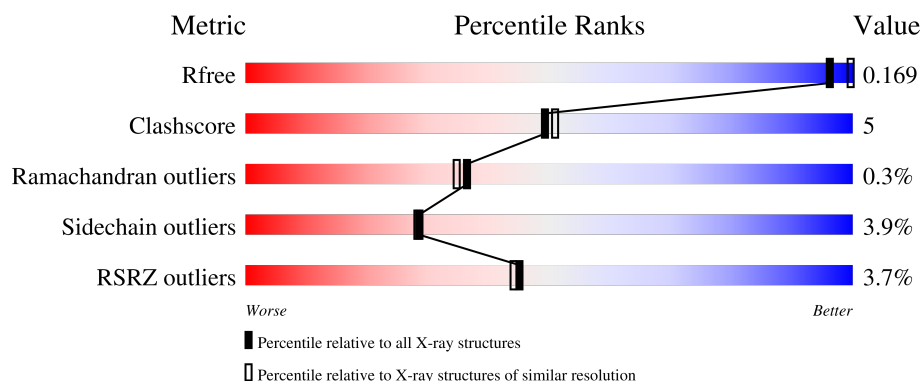
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

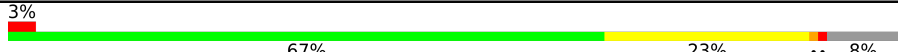
The reported resolution of this entry is 2.00 Å.

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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	359	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2887 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 Cellulase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	Se	0	0	0
			2707	1747	467	489	4			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	MSE	-	expression tag	UNP I6PLH5
A	6	GLY	-	expression tag	UNP I6PLH5
A	7	SER	-	expression tag	UNP I6PLH5
A	8	SER	-	expression tag	UNP I6PLH5
A	9	HIS	-	expression tag	UNP I6PLH5
A	10	HIS	-	expression tag	UNP I6PLH5
A	11	HIS	-	expression tag	UNP I6PLH5
A	12	HIS	-	expression tag	UNP I6PLH5
A	13	HIS	-	expression tag	UNP I6PLH5
A	14	HIS	-	expression tag	UNP I6PLH5
A	15	SER	-	expression tag	UNP I6PLH5
A	16	SER	-	expression tag	UNP I6PLH5
A	17	GLY	-	expression tag	UNP I6PLH5
A	18	LEU	-	expression tag	UNP I6PLH5
A	19	VAL	-	expression tag	UNP I6PLH5
A	20	PRO	-	expression tag	UNP I6PLH5
A	21	ARG	-	expression tag	UNP I6PLH5
A	22	GLY	-	expression tag	UNP I6PLH5
A	23	SER	-	expression tag	UNP I6PLH5
A	24	HIS	-	expression tag	UNP I6PLH5
A	25	MSE	-	expression tag	UNP I6PLH5

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

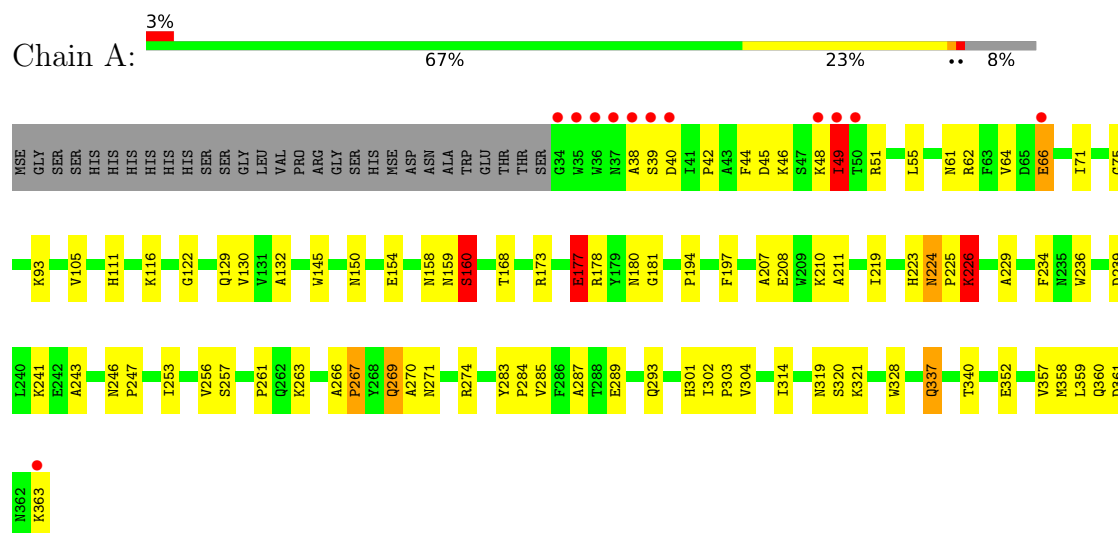
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	150	Total	O	0	0
			150	150		

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: Cellulase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	137.77 Å 137.77 Å 137.77 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.12 – 2.00 28.12 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (28.12-2.00) 99.7 (28.12-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.64 (at 1.99 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.154 , 0.169 0.155 , 0.169	Depositor DCC
R_{free} test set	2984 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	37.2	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.025 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2887	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% 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	2.09	57/2789 (2.0%)	1.67	37/3786 (1.0%)

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	A	0	1

All (57) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	358	MSE	SE-CE	-16.82	1.45	1.95
1	A	160	SER	C-O	10.22	1.37	1.24
1	A	267	PRO	N-CA	9.30	1.63	1.47
1	A	357	VAL	CA-CB	8.58	1.65	1.54
1	A	160	SER	CA-C	7.88	1.63	1.52
1	A	207	ALA	CA-CB	7.80	1.65	1.53
1	A	211	ALA	CA-CB	7.68	1.65	1.53
1	A	319	ASN	N-CA	6.91	1.54	1.46
1	A	266	ALA	CA-C	6.86	1.61	1.53
1	A	40	ASP	CA-CB	6.84	1.64	1.53
1	A	270	ALA	CA-C	6.78	1.61	1.52
1	A	243	ALA	CA-CB	6.59	1.63	1.53
1	A	226	LYS	CE-NZ	6.45	1.68	1.49
1	A	197	PHE	N-CA	6.42	1.54	1.46
1	A	289	GLU	C-O	6.21	1.31	1.23
1	A	229	ALA	CA-CB	6.14	1.63	1.53
1	A	49	ILE	CA-CB	-6.12	1.47	1.54
1	A	256	VAL	CA-CB	5.90	1.63	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	359	LEU	N-CA	5.89	1.53	1.46
1	A	130	VAL	CA-CB	5.86	1.61	1.54
1	A	234	PHE	CA-CB	5.83	1.61	1.53
1	A	159	ASN	CA-C	-5.81	1.44	1.52
1	A	197	PHE	CA-C	5.79	1.60	1.52
1	A	352	GLU	CG-CD	5.78	1.66	1.52
1	A	271	ASN	N-CA	5.77	1.53	1.46
1	A	284	PRO	CA-CB	5.75	1.61	1.53
1	A	64	VAL	CA-CB	5.73	1.65	1.55
1	A	150	ASN	C-O	5.66	1.30	1.24
1	A	253	ILE	CB-CG1	-5.59	1.42	1.53
1	A	132	ALA	CA-CB	5.58	1.62	1.53
1	A	321	LYS	C-O	-5.56	1.16	1.24
1	A	253	ILE	CB-CG2	5.54	1.70	1.52
1	A	226	LYS	CB-CG	5.53	1.69	1.52
1	A	178	ARG	CB-CG	-5.53	1.35	1.52
1	A	168	THR	N-CA	5.48	1.52	1.46
1	A	145	TRP	N-CA	5.47	1.53	1.46
1	A	181	GLY	CA-C	5.47	1.59	1.51
1	A	122	GLY	N-CA	5.45	1.52	1.45
1	A	270	ALA	CA-CB	5.40	1.61	1.53
1	A	285	VAL	C-O	5.39	1.29	1.24
1	A	208	GLU	CG-CD	5.39	1.65	1.52
1	A	194	PRO	CA-CB	5.38	1.60	1.53
1	A	274	ARG	CG-CD	5.38	1.68	1.52
1	A	363	LYS	C-O	5.33	1.34	1.23
1	A	45	ASP	CA-C	5.30	1.59	1.52
1	A	328	TRP	CA-C	5.28	1.59	1.52
1	A	75	GLY	N-CA	5.26	1.50	1.45
1	A	224	ASN	N-CA	5.21	1.52	1.46
1	A	105	VAL	CA-CB	5.19	1.62	1.55
1	A	129	GLN	CD-OE1	5.16	1.33	1.23
1	A	42	PRO	CA-C	5.13	1.58	1.52
1	A	116	LYS	N-CA	5.13	1.52	1.46
1	A	340	THR	CA-CB	5.08	1.61	1.53
1	A	320	SER	CA-CB	5.05	1.61	1.53
1	A	225	PRO	C-O	-5.04	1.17	1.24
1	A	154	GLU	CG-CD	5.03	1.64	1.52
1	A	239	ASP	CA-CB	5.01	1.60	1.53

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	62	ARG	NE-CZ-NH2	-10.12	110.10	119.20
1	A	62	ARG	NE-CZ-NH1	9.55	131.05	121.50
1	A	160	SER	N-CA-C	8.53	122.78	112.38
1	A	180	ASN	CA-C-N	-8.51	110.04	122.46
1	A	180	ASN	C-N-CA	-8.51	110.04	122.46
1	A	253	ILE	CB-CG1-CD1	-8.00	97.00	113.80
1	A	358	MSE	CG-SE-CE	7.77	116.02	98.92
1	A	226	LYS	CB-CG-CD	7.38	128.28	111.30
1	A	178	ARG	CG-CD-NE	-7.38	95.77	112.00
1	A	71	ILE	CB-CA-C	-6.92	100.31	110.62
1	A	40	ASP	CA-C-N	-6.86	117.65	122.59
1	A	40	ASP	C-N-CA	-6.86	117.65	122.59
1	A	49	ILE	N-CA-CB	-6.54	102.39	111.41
1	A	39	SER	N-CA-C	6.49	116.81	108.24
1	A	160	SER	CA-C-O	6.32	127.13	119.49
1	A	266	ALA	C-N-CD	-6.31	106.72	120.60
1	A	266	ALA	CA-C-N	6.30	142.13	127.00
1	A	266	ALA	C-N-CA	6.30	142.13	127.00
1	A	197	PHE	N-CA-CB	-6.27	99.89	110.49
1	A	49	ILE	CA-CB-CG2	6.18	121.01	110.50
1	A	302	ILE	N-CA-C	-5.99	103.45	109.02
1	A	130	VAL	CB-CA-C	-5.79	104.56	111.97
1	A	239	ASP	CB-CA-C	-5.78	100.41	109.84
1	A	62	ARG	CD-NE-CZ	5.64	132.30	124.40
1	A	177	GLU	N-CA-CB	-5.48	102.05	110.16
1	A	93	LYS	N-CA-C	5.47	119.05	112.38
1	A	49	ILE	CB-CA-C	5.38	118.59	110.63
1	A	177	GLU	CB-CG-CD	-5.37	103.47	112.60
1	A	358	MSE	CA-CB-CG	-5.29	103.52	114.10
1	A	283	TYR	CA-C-N	-5.28	114.52	119.85
1	A	283	TYR	C-N-CA	-5.28	114.52	119.85
1	A	234	PHE	CA-CB-CG	-5.28	108.53	113.80
1	A	359	LEU	N-CA-C	-5.23	105.76	111.82
1	A	261	PRO	CB-CA-C	-5.21	104.08	112.21
1	A	51	ARG	NE-CZ-NH2	-5.13	114.58	119.20
1	A	61	ASN	CA-C-N	-5.04	114.68	122.59
1	A	61	ASN	C-N-CA	-5.04	114.68	122.59

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	38	ALA	Peptide

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	2707	0	2595	27	0
2	A	30	0	40	7	0
3	A	150	0	0	3	0
All	All	2887	0	2635	29	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 5.

All (29) 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:226:LYS:CE	1:A:226:LYS:NZ	1.68	1.51
1:A:269:GLN:HE21	1:A:269:GLN:H	1.25	0.84
1:A:66:GLU:H	1:A:66:GLU:CD	1.91	0.78
2:A:603:GOL:H31	3:A:820:HOH:O	1.84	0.77
1:A:224:ASN:HD21	1:A:226:LYS:NZ	1.85	0.74
1:A:223:HIS:HE1	3:A:715:HOH:O	1.70	0.74
1:A:224:ASN:HD21	1:A:226:LYS:HZ2	1.43	0.65
1:A:66:GLU:CD	1:A:66:GLU:N	2.58	0.62
1:A:219:ILE:O	1:A:223:HIS:HD2	1.83	0.61
1:A:337:GLN:NE2	1:A:337:GLN:H	2.00	0.59
1:A:111:HIS:NE2	2:A:601:GOL:H31	2.21	0.56
1:A:263:LYS:NZ	2:A:604:GOL:O1	2.32	0.56
1:A:269:GLN:H	1:A:269:GLN:NE2	2.01	0.55
1:A:360:GLN:HG2	1:A:361:ASP:OD1	2.08	0.52
1:A:210:LYS:HE3	1:A:246:ASN:HD21	1.75	0.51
1:A:224:ASN:ND2	1:A:226:LYS:NZ	2.57	0.48
1:A:303:PRO:HG3	2:A:604:GOL:H12	1.99	0.45
1:A:236:TRP:CD2	2:A:604:GOL:H2	2.52	0.45
1:A:301:HIS:O	1:A:304:VAL:HG22	2.17	0.45
1:A:257:SER:O	1:A:287:ALA:HA	2.17	0.44
1:A:210:LYS:HE2	1:A:247:PRO:O	2.17	0.44
1:A:210:LYS:HE3	1:A:246:ASN:ND2	2.33	0.44
1:A:293:GLN:HE22	1:A:337:GLN:HE22	1.64	0.43
1:A:111:HIS:NE2	2:A:601:GOL:C3	2.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:PHE:CZ	1:A:49:ILE:HD12	2.54	0.42
1:A:173:ARG:O	1:A:177:GLU:HB3	2.20	0.41
2:A:603:GOL:C3	3:A:820:HOH:O	2.55	0.41
1:A:55:LEU:HD11	1:A:226:LYS:HB2	2.03	0.41
1:A:158:ASN:OD1	1:A:160:SER:HB3	2.21	0.41

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	328/359 (91%)	321 (98%)	6 (2%)	1 (0%)	36	35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	267	PRO

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	281/301 (93%)	270 (96%)	11 (4%)	28	28

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LYS
1	A	48	LYS
1	A	49	ILE
1	A	66	GLU
1	A	160	SER
1	A	177	GLU
1	A	226	LYS
1	A	241	LYS
1	A	269	GLN
1	A	314	ILE
1	A	337	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	135	ASN
1	A	213	ASN
1	A	223	HIS
1	A	224	ASN
1	A	246	ASN
1	A	269	GLN
1	A	337	GLN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

5 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	A	601	-	5,5,5	0.41	0	5,5,5	1.47	1 (20%)
2	GOL	A	603	-	5,5,5	0.90	0	5,5,5	1.16	0
2	GOL	A	602	-	5,5,5	0.66	0	5,5,5	1.27	1 (20%)
2	GOL	A	604	-	5,5,5	0.82	0	5,5,5	1.46	1 (20%)
2	GOL	A	605	-	5,5,5	0.43	0	5,5,5	1.02	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	A	601	-	-	2/4/4/4	-
2	GOL	A	603	-	-	2/4/4/4	-
2	GOL	A	602	-	-	2/4/4/4	-
2	GOL	A	604	-	-	4/4/4/4	-
2	GOL	A	605	-	-	2/4/4/4	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	604	GOL	C3-C2-C1	2.92	122.52	111.80
2	A	601	GOL	O1-C1-C2	-2.87	97.44	110.38
2	A	602	GOL	O2-C2-C1	-2.14	100.32	109.18

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
2	A	602	GOL	C1-C2-C3-O3
2	A	603	GOL	C1-C2-C3-O3
2	A	603	GOL	O2-C2-C3-O3
2	A	605	GOL	C1-C2-C3-O3
2	A	604	GOL	O1-C1-C2-C3
2	A	601	GOL	O2-C2-C3-O3
2	A	605	GOL	O2-C2-C3-O3
2	A	602	GOL	O2-C2-C3-O3
2	A	604	GOL	C1-C2-C3-O3
2	A	604	GOL	O1-C1-C2-O2
2	A	604	GOL	O2-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	GOL	2	0
2	A	603	GOL	2	0
2	A	604	GOL	3	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	326/359 (90%)	-0.10	12 (3%) 45 44	29, 41, 66, 88	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	40	ASP	5.9
1	A	36	TRP	5.8
1	A	34	GLY	4.8
1	A	66	GLU	4.7
1	A	363	LYS	4.0
1	A	49	ILE	3.4
1	A	37	ASN	3.3
1	A	39	SER	2.9
1	A	48	LYS	2.6
1	A	38	ALA	2.6
1	A	35	TRP	2.5
1	A	50	THR	2.1

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(Å ²)	Q<0.9
2	GOL	A	604	6/6	0.80	0.24	58,70,77,82	3
2	GOL	A	601	6/6	0.86	0.17	64,74,76,77	0
2	GOL	A	605	6/6	0.86	0.22	77,83,83,86	0
2	GOL	A	602	6/6	0.91	0.15	56,66,67,69	0
2	GOL	A	603	6/6	0.94	0.14	36,59,71,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.