



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 11:16 AM GMT

PDB ID : 3OGV
Title : Complex structure of beta-galactosidase from *Trichoderma reesei* with PETG
Authors : Maksimainen, M.; Rouvinen, J.
Deposited on : 2010-08-17
Resolution : 1.40 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

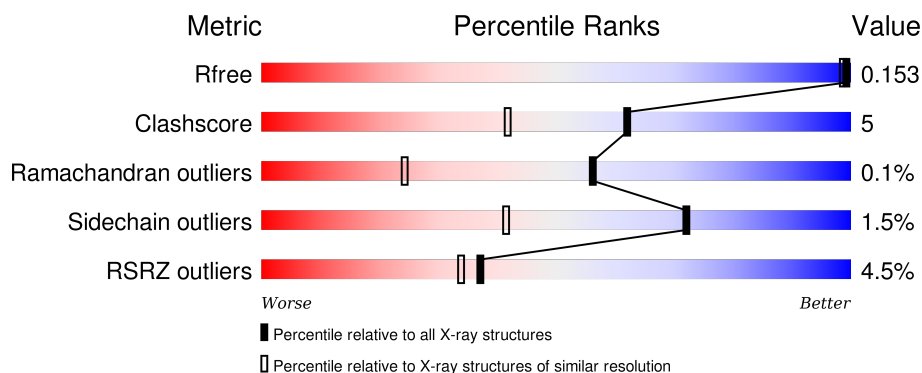
1 Overall quality at a glance

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 ≥ 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	1003	<div> <div>4%</div> <div>82%</div> <div>14%</div> <div>••</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	MAN	A	1031	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	NAG	A	1040	-	-	-	X
5	NAG	A	1041	-	-	-	X
6	NAG	A	1043	-	-	-	X

2 Entry composition [i](#)

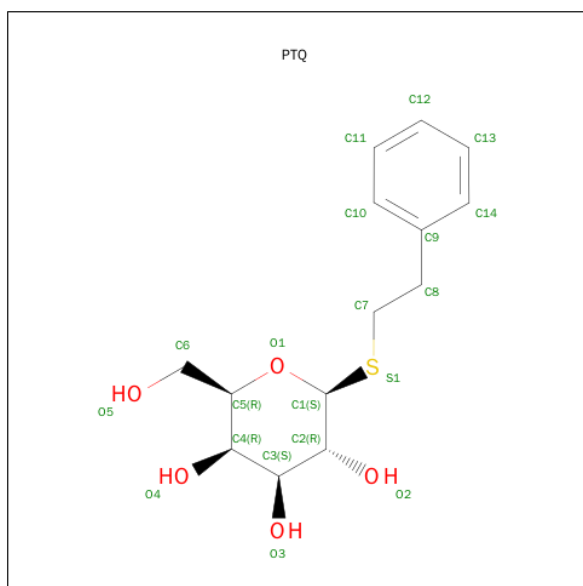
There are 7 unique types of molecules in this entry. The entry contains 9065 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Beta-galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	986	Total	C	N	O	S	0	10	0
			7674	4937	1292	1437	8			

- Molecule 2 is SUGAR (2-PHENYLETHYL 1-THIO-BETA-D-GALACTOPYRANOSIDE) (three-letter code: PTQ) (formula: C₁₄H₂₀O₅S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	S	0	0
			20	14	5	1		

- Molecule 3 is a polymer of unknown type called SUGAR (7-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	7	Total	C	N	O	0	0
			83	46	2	35		

- Molecule 4 is a polymer of unknown type called SUGAR (8-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	8	Total	C	N	O	0	0
			94	52	2	40		

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	2	Total	C	N	O	0	0
			28	16	2	10		

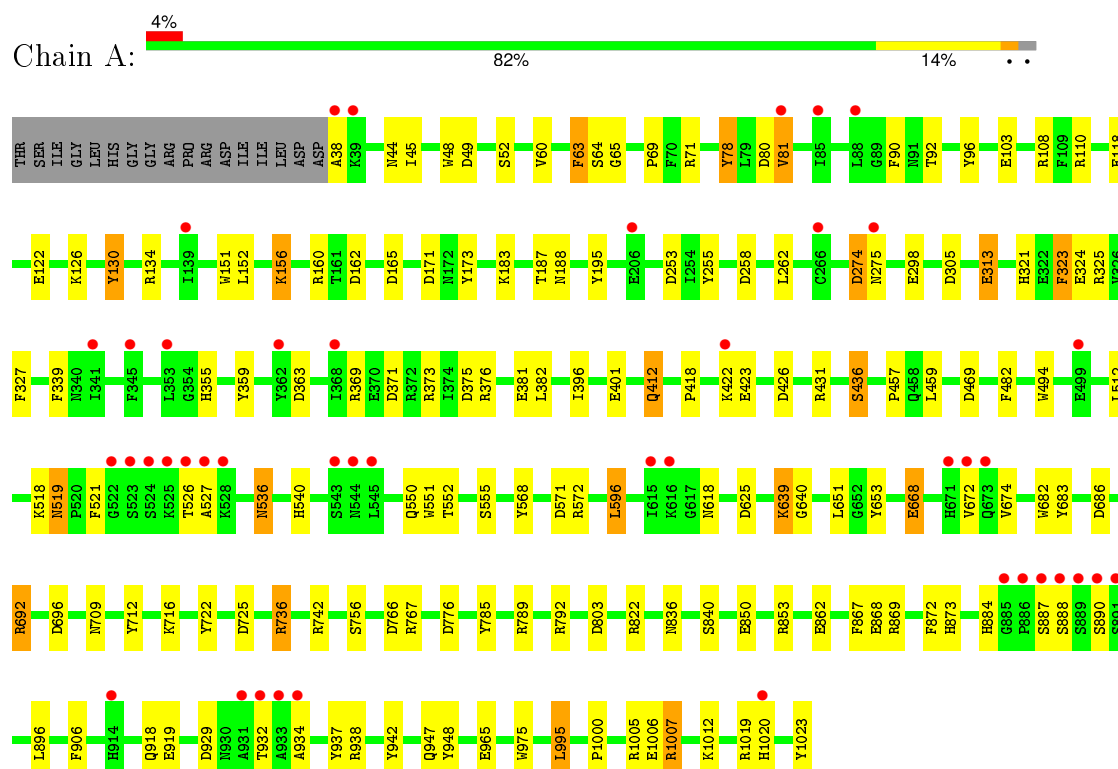
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1138	Total	O	0	0
			1138	1138		

3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: Beta-galactosidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	67.60 Å 68.70 Å 81.70 Å 108.50° 97.70° 114.50°	Depositor
Resolution (Å)	42.76 – 1.40 42.76 – 1.35	Depositor EDS
% Data completeness (in resolution range)	95.0 (42.76-1.40) 77.7 (42.76-1.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 1.35 Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.132 , 0.166 0.145 , 0.153	Depositor DCC
R_{free} test set	11034 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	10.9	Xtriage
Anisotropy	0.483	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 59.5	EDS
Estimated twinning fraction	0.015 for k,h,-h-k-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 244951 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9065	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: GLC, BMA, NAG, PTQ, MAN

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	1.49	46/7915 (0.6%)	1.38	87/10787 (0.8%)

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

The worst 5 of 46 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	313	GLU	CD-OE1	11.05	1.37	1.25
1	A	551	TRP	CB-CG	8.83	1.66	1.50
1	A	324	GLU	CD-OE1	8.36	1.34	1.25
1	A	436	SER	CA-CB	8.24	1.65	1.52
1	A	325	ARG	CG-CD	7.67	1.71	1.51

The worst 5 of 87 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	572	ARG	NE-CZ-NH2	-13.98	113.31	120.30
1	A	938	ARG	NE-CZ-NH2	-13.69	113.45	120.30
1	A	938	ARG	NE-CZ-NH1	12.32	126.46	120.30
1	A	431	ARG	NE-CZ-NH1	11.73	126.16	120.30
1	A	162	ASP	CB-CG-OD2	-11.16	108.26	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	736	ARG	Sidechain

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	7674	0	7490	83	2
2	A	20	0	20	1	0
3	A	83	0	70	2	0
4	A	94	0	79	0	0
5	A	28	0	26	0	0
6	A	28	0	25	1	0
7	A	1138	0	0	44	3
All	All	9065	0	7710	84	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 5.

The worst 5 of 84 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:108:ARG:HD2	7:A:2107:HOH:O	1.34	1.22
1:A:1012:LYS:HE3	7:A:1986:HOH:O	1.42	1.19
1:A:1020:HIS:CD2	7:A:2077:HOH:O	2.04	1.11
1:A:69:PRO:HG2	7:A:1597:HOH:O	1.50	1.10
1:A:1006:GLU:HG3	7:A:1815:HOH:O	1.51	1.07

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)
7:A:1305:HOH:O	7:A:2204:HOH:O[1_556]	1.72	0.48
1:A:540:HIS:CE1	1:A:1020:HIS:CE1[1_544]	1.88	0.32
7:A:1300:HOH:O	7:A:2160:HOH:O[1_455]	2.10	0.10
1:A:540:HIS:CE1	1:A:1020:HIS:NE2[1_544]	2.12	0.08
7:A:1873:HOH:O	7:A:1988:HOH:O[1_655]	2.12	0.08

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	994/1003 (99%)	965 (97%)	28 (3%)	1 (0%)	56 24

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	519	ASN

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	815/819 (100%)	803 (98%)	12 (2%)	72 41

5 of 12 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	618	ASN
1	A	639	LYS
1	A	947	GLN
1	A	536	ASN
1	A	932	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	618	ASN
1	A	673	GLN
1	A	884	HIS
1	A	536	ASN
1	A	873	HIS

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 ⓘ

17 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	$\# Z > 2$
3	NAG	A	1025	1,3	14,14,15	0.50	0	15,19,21	1.79	3 (20%)
3	NAG	A	1026	3	14,14,15	0.72	0	15,19,21	1.15	2 (13%)
3	BMA	A	1027	3	11,11,12	1.18	1 (9%)	14,15,17	1.02	1 (7%)
3	MAN	A	1028	3	11,11,12	0.83	0	14,15,17	1.12	1 (7%)
3	MAN	A	1029	3	11,11,12	0.95	0	14,15,17	1.65	1 (7%)
3	MAN	A	1030	3	11,11,12	1.24	2 (18%)	14,15,17	1.44	2 (14%)
3	MAN	A	1031	3	11,11,12	1.08	0	14,15,17	2.39	7 (50%)
4	NAG	A	1032	1,4	14,14,15	1.88	4 (28%)	15,19,21	1.65	2 (13%)
4	NAG	A	1033	4	14,14,15	0.75	0	15,19,21	1.26	1 (6%)
4	BMA	A	1034	4	11,11,12	1.01	1 (9%)	14,15,17	1.31	2 (14%)
4	MAN	A	1035	4	11,11,12	1.12	1 (9%)	14,15,17	2.25	3 (21%)
4	MAN	A	1036	4	11,11,12	0.93	1 (9%)	14,15,17	1.84	6 (42%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MAN	A	1037	4	11,11,12	1.02	0	14,15,17	1.79	3 (21%)
4	MAN	A	1038	4	11,11,12	0.78	0	14,15,17	1.32	3 (21%)
4	GLC	A	1039	4	11,11,12	0.78	0	14,15,17	2.24	4 (28%)
6	NAG	A	1042	1,6	14,14,15	1.04	1 (7%)	15,19,21	2.00	6 (40%)
6	NAG	A	1043	6	14,14,15	0.56	0	15,19,21	1.61	3 (20%)

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	NAG	A	1025	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	1026	3	-	0/6/23/26	0/1/1/1
3	BMA	A	1027	3	-	0/2/19/22	0/1/1/1
3	MAN	A	1028	3	-	0/2/19/22	0/1/1/1
3	MAN	A	1029	3	-	0/2/19/22	0/1/1/1
3	MAN	A	1030	3	-	0/2/19/22	0/1/1/1
3	MAN	A	1031	3	-	0/2/19/22	0/1/1/1
4	NAG	A	1032	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	1033	4	-	0/6/23/26	0/1/1/1
4	BMA	A	1034	4	-	0/2/19/22	0/1/1/1
4	MAN	A	1035	4	-	0/2/19/22	0/1/1/1
4	MAN	A	1036	4	-	0/2/19/22	0/1/1/1
4	MAN	A	1037	4	-	0/2/19/22	0/1/1/1
4	MAN	A	1038	4	-	0/2/19/22	0/1/1/1
4	GLC	A	1039	4	-	0/2/19/22	0/1/1/1
6	NAG	A	1042	1,6	-	0/6/23/26	0/1/1/1
6	NAG	A	1043	6	-	0/6/23/26	0/1/1/1

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1032	NAG	O5-C1	-3.81	1.37	1.43
6	A	1042	NAG	C1-C2	-2.80	1.48	1.52
4	A	1035	MAN	O5-C1	-2.75	1.39	1.43
4	A	1032	NAG	C2-N2	-2.32	1.42	1.46
3	A	1030	MAN	O2-C2	-2.23	1.38	1.43

The worst 5 of 50 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1031	MAN	C6-C5-C4	-5.43	99.62	113.02
6	A	1043	NAG	C3-C4-C5	-4.06	103.12	110.20
3	A	1025	NAG	C2-N2-C7	-3.76	118.21	123.04
4	A	1033	NAG	O4-C4-C3	-3.69	102.03	110.34
4	A	1037	MAN	C1-C2-C3	-3.61	105.27	109.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1028	MAN	1	0
3	A	1029	MAN	1	0
6	A	1042	NAG	1	0

5.6 Ligand geometry

3 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	PTQ	A	1024	-	21,21,21	1.46	4 (19%)	26,28,28	1.49	5 (19%)
5	NAG	A	1040	1	14,14,15	1.27	1 (7%)	15,19,21	1.68	3 (20%)
5	NAG	A	1041	1	14,14,15	0.82	0	15,19,21	2.09	5 (33%)

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	PTQ	A	1024	-	-	0/8/28/28	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	1040	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1041	1	-	0/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1024	PTQ	C13-C14	-2.34	1.34	1.38
2	A	1024	PTQ	C1-C2	-2.02	1.49	1.53
2	A	1024	PTQ	C1-S1	2.48	1.84	1.80
2	A	1024	PTQ	O1-C5	2.88	1.51	1.44
5	A	1040	NAG	O7-C7	3.98	1.32	1.23

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1041	NAG	O7-C7-C8	-4.43	113.93	122.06
5	A	1040	NAG	O6-C6-C5	-3.30	100.43	111.33
5	A	1040	NAG	C8-C7-N2	-3.17	110.05	116.11
5	A	1041	NAG	O4-C4-C5	-2.97	101.36	109.24
5	A	1041	NAG	C2-N2-C7	-2.74	119.52	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1024	PTQ	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	986/1003 (98%)	0.16	44 (4%) 37 34	6, 12, 26, 61	0

The worst 5 of 44 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	526	THR	11.8
1	A	933	ALA	10.1
1	A	932	THR	7.6
1	A	523	SER	7.0
1	A	525	LYS	6.8

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, 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	LLDF	B-factors(Å ²)	Q<0.9
6	NAG	A	1043	14/15	0.70	0.34	14.94	35,40,51,51	0
3	MAN	A	1031	11/12	0.93	0.14	2.41	15,16,22,25	0
6	NAG	A	1042	14/15	0.88	0.15	1.09	22,30,38,41	0
4	MAN	A	1037	11/12	0.96	0.10	0.77	12,13,16,16	0
4	NAG	A	1032	14/15	0.96	0.08	-0.23	11,13,22,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	GLC	A	1039	11/12	0.95	0.09	-0.23	13,15,19,21	0
3	MAN	A	1029	11/12	0.97	0.07	-0.37	10,12,19,29	0
3	NAG	A	1025	14/15	0.97	0.07	-1.02	8,11,18,21	0
3	NAG	A	1026	14/15	0.98	0.06	-1.57	8,10,15,16	0
4	NAG	A	1033	14/15	0.96	0.10	-	10,14,32,34	0
3	MAN	A	1028	11/12	0.96	0.07	-	9,11,16,22	0
4	MAN	A	1036	11/12	0.94	0.11	-	14,18,23,31	0
3	MAN	A	1030	11/12	0.89	0.27	-	21,25,36,37	0
4	MAN	A	1038	11/12	0.97	0.07	-	13,13,15,16	0
3	BMA	A	1027	11/12	0.98	0.06	-	9,11,14,15	0
4	BMA	A	1034	11/12	0.97	0.09	-	12,13,17,18	0
4	MAN	A	1035	11/12	0.82	0.24	-	27,31,37,40	0

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. 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, 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	LLDF	B-factors(\AA^2)	Q<0.9
5	NAG	A	1041	14/15	0.90	0.20	5.73	24,28,35,41	0
5	NAG	A	1040	14/15	0.79	0.23	3.52	28,36,45,48	0
2	PTQ	A	1024	20/20	0.97	0.09	-0.34	6,8,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.