



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 2, 2016 – 01:15 AM EDT

PDB ID : 5JP0
Title : Bacteroides ovatus Xyloglucan PUL GH3B with bound glucose
Authors : Hemsworth, G.R.; Thompson, A.J.; Stepper, J.; Sobala, L.F.; Coyle, T.; Larsbrink, J.; Spadiut, O.; Stubbs, K.A.; Brumer, H.; Davies, G.J.
Deposited on : 2016-05-03
Resolution : 2.30 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

<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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027939
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) : rb-20027939

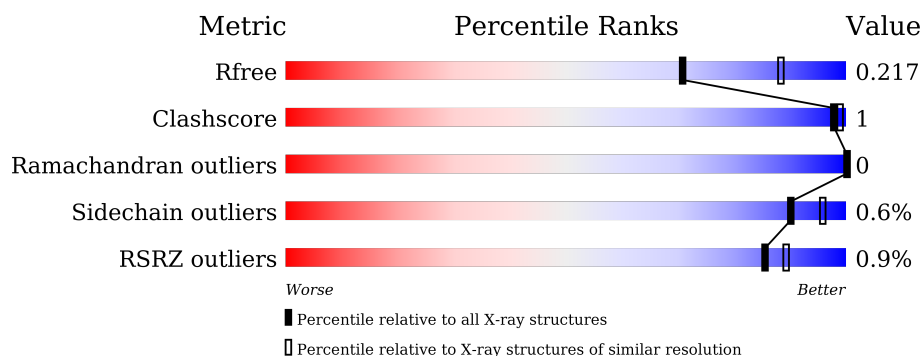
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	772	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> % 93% . . </div> </div>
1	B	772	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; padding: 0 5px;"> % 93% . . </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11996 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.

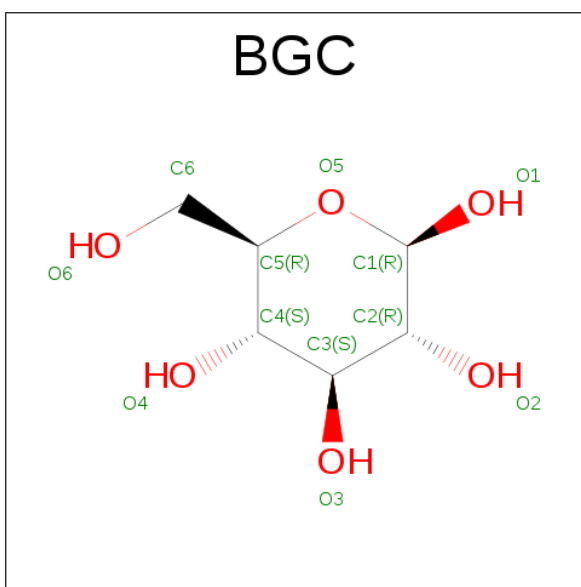
- Molecule 1 is a protein called Beta-glucosidase BoGH3B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	745	Total	C	N	O	S	0	0	0
			5757	3660	964	1098	35			
1	B	745	Total	C	N	O	S	0	0	0
			5751	3657	961	1098	35			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	787	LEU	-	expression tag	UNP A7LXU3
A	788	GLU	-	expression tag	UNP A7LXU3
A	789	HIS	-	expression tag	UNP A7LXU3
A	790	HIS	-	expression tag	UNP A7LXU3
A	791	HIS	-	expression tag	UNP A7LXU3
A	792	HIS	-	expression tag	UNP A7LXU3
A	793	HIS	-	expression tag	UNP A7LXU3
A	794	HIS	-	expression tag	UNP A7LXU3
B	787	LEU	-	expression tag	UNP A7LXU3
B	788	GLU	-	expression tag	UNP A7LXU3
B	789	HIS	-	expression tag	UNP A7LXU3
B	790	HIS	-	expression tag	UNP A7LXU3
B	791	HIS	-	expression tag	UNP A7LXU3
B	792	HIS	-	expression tag	UNP A7LXU3
B	793	HIS	-	expression tag	UNP A7LXU3
B	794	HIS	-	expression tag	UNP A7LXU3

- Molecule 2 is BETA-D-GLUCOSE (three-letter code: BGC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		
2	B	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		

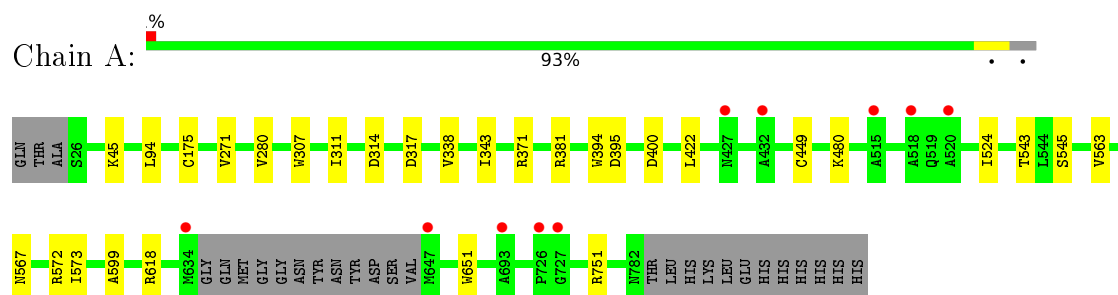
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	210	Total	O	0	0
			210	210		
4	B	252	Total	O	0	0
			252	252		

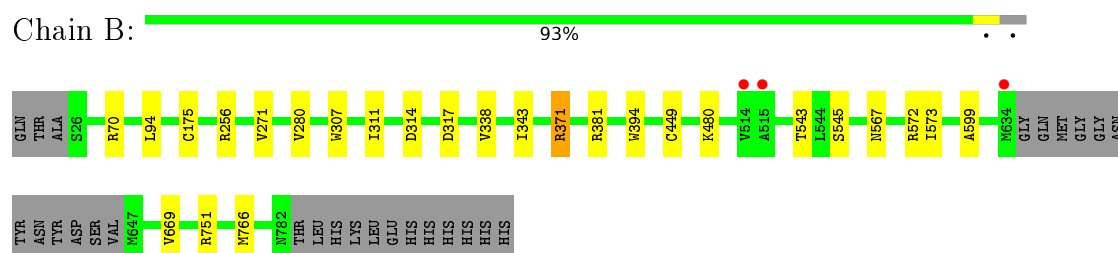
3 Residue-property plots [i](#)

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-glucosidase BoGH3B



- Molecule 1: Beta-glucosidase BoGH3B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	97.17Å 104.07Å 160.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.57 – 2.30 47.57 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (47.57-2.30) 99.9 (47.57-2.30)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.177 , 0.211 0.185 , 0.217	Depositor DCC
R_{free} test set	3561 reflections (5.14%)	DCC
Wilson B-factor (Å ²)	30.8	Xtriage
Anisotropy	0.769	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 27.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11996	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.43% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, BGC

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	0.53	0/5885	0.74	11/8002 (0.1%)
1	B	0.54	0/5879	0.83	9/7995 (0.1%)
All	All	0.53	0/11764	0.79	20/15997 (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	B	0	1

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	371	ARG	NE-CZ-NH2	-31.25	104.68	120.30
1	B	371	ARG	NE-CZ-NH1	26.42	133.51	120.30
1	A	371	ARG	NE-CZ-NH2	18.82	129.71	120.30
1	A	371	ARG	NE-CZ-NH1	-15.19	112.70	120.30
1	B	371	ARG	CD-NE-CZ	9.51	136.92	123.60
1	A	395	ASP	CB-CG-OD1	6.50	124.15	118.30
1	A	618	ARG	NE-CZ-NH1	6.05	123.33	120.30
1	A	572	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	A	395	ASP	CB-CG-OD2	-5.71	113.16	118.30
1	B	751	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	A	572	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	B	572	ARG	NE-CZ-NH2	-5.43	117.59	120.30
1	B	371	ARG	CG-CD-NE	-5.41	100.43	111.80
1	A	381	ARG	NE-CZ-NH1	5.24	122.92	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	751	ARG	NE-CZ-NH1	5.16	122.88	120.30
1	A	400	ASP	CB-CG-OD1	5.12	122.91	118.30
1	B	572	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	B	381	ARG	NE-CZ-NH1	5.11	122.86	120.30
1	A	618	ARG	NE-CZ-NH2	-5.05	117.77	120.30
1	B	751	ARG	NE-CZ-NH2	-5.05	117.78	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	371	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	5757	0	5588	14	0
1	B	5751	0	5577	12	0
2	A	12	0	12	1	0
2	B	12	0	12	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	210	0	0	0	0
4	B	252	0	0	0	0
All	All	11996	0	11189	26	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 1.

All (26) 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:B:314:ASP:OD1	2:B:801:BGC:H1	1.76	0.84
1:A:311:ILE:HG22	1:A:343:ILE:HD13	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:314:ASP:OD1	2:A:801:BGC:H1	2.01	0.60
1:B:311:ILE:HG22	1:B:343:ILE:HD13	1.85	0.58
1:A:524:ILE:HG22	1:A:563:VAL:HB	1.88	0.54
1:A:175:CYS:HA	1:A:394:TRP:CZ2	2.46	0.51
1:B:175:CYS:HA	1:B:394:TRP:CZ2	2.47	0.50
1:B:271:VAL:HG13	1:B:307:TRP:CD1	2.47	0.48
1:A:338:VAL:CG1	1:A:343:ILE:HG13	2.43	0.48
1:A:449:CYS:HB3	1:A:567:ASN:OD1	2.14	0.48
1:B:338:VAL:CG1	1:B:343:ILE:HG13	2.45	0.47
1:B:449:CYS:HB3	1:B:567:ASN:OD1	2.14	0.47
1:A:280:VAL:HG21	1:A:343:ILE:HD11	1.98	0.45
1:A:480:LYS:HG3	1:A:599:ALA:HB1	1.98	0.45
1:B:543:THR:HA	1:B:573:ILE:HB	2.00	0.44
1:B:480:LYS:HG3	1:B:599:ALA:HB1	1.98	0.44
1:B:280:VAL:HG21	1:B:343:ILE:HD11	1.98	0.44
1:B:669:VAL:HG12	1:B:766:MET:HE1	2.00	0.43
1:A:271:VAL:HG13	1:A:307:TRP:CD1	2.53	0.43
1:A:543:THR:HA	1:A:573:ILE:HB	1.99	0.43
1:B:314:ASP:HB3	1:B:317:ASP:OD2	2.20	0.42
1:A:314:ASP:HB3	1:A:317:ASP:OD2	2.19	0.42
1:A:422:LEU:HD13	1:A:651:TRP:CE3	2.55	0.42
1:A:338:VAL:HG13	1:A:343:ILE:HG13	2.01	0.41
1:A:311:ILE:CG2	1:A:343:ILE:HD13	2.49	0.41
1:B:338:VAL:HG13	1:B:343:ILE:HG13	2.02	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	741/772 (96%)	717 (97%)	24 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	741/772 (96%)	717 (97%)	24 (3%)	0	100	100
All	All	1482/1544 (96%)	1434 (97%)	48 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	606/653 (93%)	603 (100%)	3 (0%)	92	97
1	B	605/653 (93%)	601 (99%)	4 (1%)	88	95
All	All	1211/1306 (93%)	1204 (99%)	7 (1%)	90	96

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	LYS
1	A	94	LEU
1	A	545	SER
1	B	70	ARG
1	B	94	LEU
1	B	256	ARG
1	B	545	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	633	ASN
1	B	633	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\# Z > 2$
2	BGC	A	801	-	12,12,12	0.62	0	17,17,17	1.72	3 (17%)
2	BGC	B	801	-	12,12,12	0.49	0	17,17,17	1.48	3 (17%)

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	BGC	A	801	-	-	0/2/22/22	0/1/1/1
2	BGC	B	801	-	-	0/2/22/22	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	BGC	C1-C2-C3	2.24	114.35	110.68
2	B	801	BGC	O5-C1-C2	2.32	114.06	110.00
2	A	801	BGC	C1-C2-C3	3.33	116.14	110.68
2	A	801	BGC	O5-C1-C2	3.53	116.18	110.00
2	B	801	BGC	C1-O5-C5	4.28	121.72	113.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	BGC	C1-O5-C5	4.54	122.23	113.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	BGC	1	0
2	B	801	BGC	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 [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	745/772 (96%)	-0.09	10 (1%) 79 84	20, 34, 58, 77	0
1	B	745/772 (96%)	-0.18	3 (0%) 93 95	20, 33, 54, 84	0
All	All	1490/1544 (96%)	-0.14	13 (0%) 85 89	20, 34, 56, 84	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	726	PRO	3.3
1	A	515	ALA	3.3
1	B	515	ALA	3.1
1	A	647	MET	2.9
1	A	693	ALA	2.8
1	A	427	ASN	2.6
1	B	514	VAL	2.6
1	A	518	ALA	2.5
1	B	634	MET	2.5
1	A	727	GLY	2.3
1	A	634	MET	2.3
1	A	520	ALA	2.2
1	A	432	ALA	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 carbohydrates 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. 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
2	BGC	A	801	12/12	0.90	0.17	1.37	24,35,45,52	0
2	BGC	B	801	12/12	0.95	0.12	-0.76	22,27,30,32	0
3	MG	A	802	1/1	0.97	0.10	-1.81	27,27,27,27	0
3	MG	B	802	1/1	0.96	0.03	-7.22	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.