



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 24, 2017 – 09:21 PM EDT

PDB ID : 5KZJ
Title : Loop Deletion mutant of Paracoccus denitrificans AztC
Authors : Yukl, E.
Deposited on : unknown
Resolution : 2.20 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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-20030345

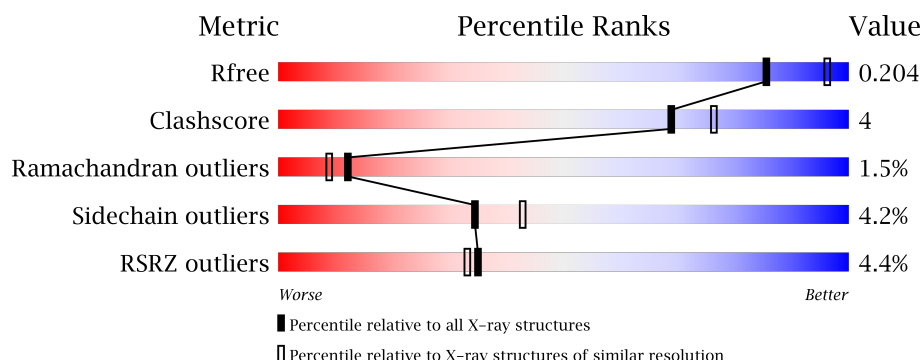
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

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	296	<div> <div>5%</div> <div> <div></div> <div>80%</div> <div>9%</div> <div>8%</div> </div> </div>
1	B	296	<div> <div>3%</div> <div> <div></div> <div>80%</div> <div>10%</div> <div>8%</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
2	ZN	A	401	-	-	-	X
2	ZN	B	401	-	-	-	X
3	GOL	A	402	-	-	-	X
3	GOL	A	403	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4194 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 Periplasmic solute binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	272	Total	C	N	O	S	0	3	0
			2013	1255	355	398	5			
1	B	271	Total	C	N	O	S	0	3	0
			2004	1251	353	395	5			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	HIS	deletion	UNP A1B2F3
A	?	-	TYR	deletion	UNP A1B2F3
A	?	-	HIS	deletion	UNP A1B2F3
A	?	-	TYR	deletion	UNP A1B2F3
A	?	-	ILE	deletion	UNP A1B2F3
A	?	-	ASP	deletion	UNP A1B2F3
A	?	-	GLY	deletion	UNP A1B2F3
A	?	-	LYS	deletion	UNP A1B2F3
A	?	-	ALA	deletion	UNP A1B2F3
A	?	-	VAL	deletion	UNP A1B2F3
A	?	-	PHE	deletion	UNP A1B2F3
A	?	-	HIS	deletion	UNP A1B2F3
A	?	-	ALA	deletion	UNP A1B2F3
B	?	-	HIS	deletion	UNP A1B2F3
B	?	-	TYR	deletion	UNP A1B2F3
B	?	-	HIS	deletion	UNP A1B2F3
B	?	-	TYR	deletion	UNP A1B2F3
B	?	-	ILE	deletion	UNP A1B2F3
B	?	-	ASP	deletion	UNP A1B2F3
B	?	-	GLY	deletion	UNP A1B2F3
B	?	-	LYS	deletion	UNP A1B2F3
B	?	-	ALA	deletion	UNP A1B2F3
B	?	-	VAL	deletion	UNP A1B2F3
B	?	-	PHE	deletion	UNP A1B2F3
B	?	-	HIS	deletion	UNP A1B2F3

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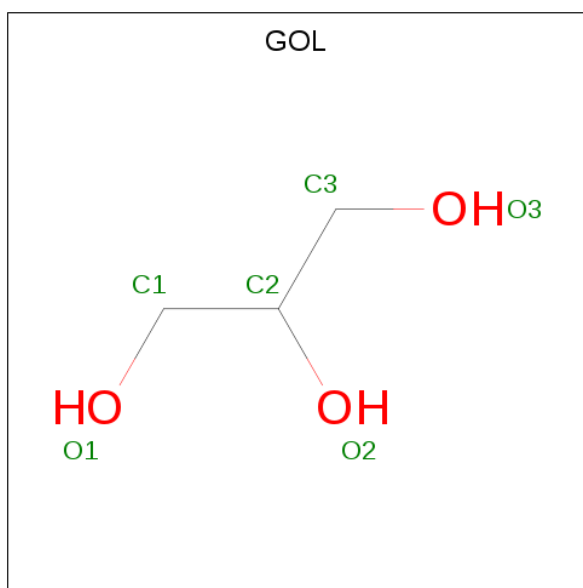
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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	ALA	deletion	UNP A1B2F3

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

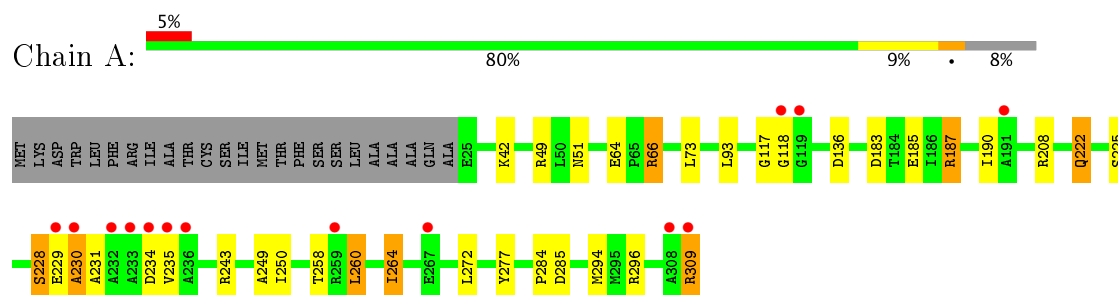
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	85	Total	O	0	0
			85	85		
4	B	77	Total	O	0	1
			78	78		

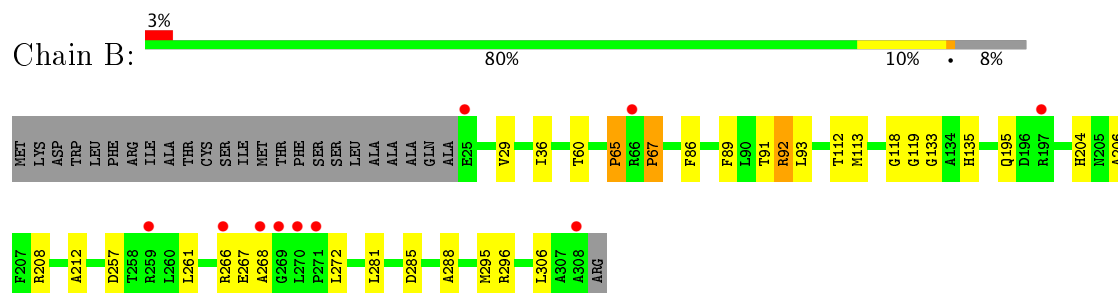
3 Residue-property plots [i](#)

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



- Molecule 1: Periplasmic solute binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.83Å 104.31Å 64.32Å 90.00° 110.70° 90.00°	Depositor
Resolution (Å)	46.47 – 2.20 46.43 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (46.47-2.20) 99.6 (46.43-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.159 , 0.200 0.167 , 0.204	Depositor DCC
R_{free} test set	1942 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	43.5	Xtriage
Anisotropy	0.208	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.043 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4194	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.59% 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: GOL, ZN

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	0.91	0/2049	1.02	6/2797 (0.2%)
1	B	0.97	0/2041	1.01	7/2786 (0.3%)
All	All	0.94	0/4090	1.01	13/5583 (0.2%)

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

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	208	ARG	NE-CZ-NH1	11.61	126.10	120.30
1	B	296	ARG	NE-CZ-NH2	-10.41	115.09	120.30
1	A	208	ARG	NE-CZ-NH2	-8.31	116.14	120.30
1	B	296	ARG	NE-CZ-NH1	7.57	124.09	120.30
1	B	208	ARG	NE-CZ-NH1	7.22	123.91	120.30
1	B	92	ARG	NE-CZ-NH2	-6.45	117.07	120.30
1	A	136	ASP	CB-CG-OD2	-6.25	112.68	118.30
1	B	92	ARG	NE-CZ-NH1	6.09	123.35	120.30
1	A	66	ARG	NE-CZ-NH2	5.78	123.19	120.30
1	A	187	ARG	NE-CZ-NH1	5.32	122.96	120.30
1	B	285	ASP	CB-CG-OD1	5.20	122.98	118.30
1	A	260	LEU	CA-CB-CG	5.16	127.16	115.30
1	B	295	MET	CG-SD-CE	-5.13	92.00	100.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	228	SER	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	2013	0	1943	18	0
1	B	2004	0	1932	13	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	12	0	16	3	0
4	A	85	0	0	5	0
4	B	78	0	0	0	0
All	All	4194	0	3891	34	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 4.

All (34) 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:185:GLU:OE1	1:A:296:ARG:NH2	2.14	0.78
1:A:260:LEU:O	1:A:264:ILE:HD13	1.93	0.68
1:B:266:ARG:O	1:B:268:ALA:N	2.27	0.67
1:B:112:THR:OG1	1:B:135[A]:HIS:CD2	2.52	0.62
3:A:403:GOL:C3	4:A:547:HOH:O	2.50	0.59
1:A:222[A]:GLN:NE2	1:A:225:SER:O	2.36	0.58
3:A:403:GOL:O3	4:A:501:HOH:O	2.16	0.58
1:A:64[A]:GLU:O	1:A:66:ARG:HD3	2.04	0.58
1:A:183:ASP:O	1:A:187:ARG:HG3	2.04	0.57
1:A:64[B]:GLU:O	1:A:66:ARG:HD3	2.06	0.55
1:B:266:ARG:C	1:B:268:ALA:H	2.09	0.55
1:A:250:ILE:HD12	1:A:250:ILE:N	2.22	0.55
1:A:230:ALA:O	1:A:231:ALA:C	2.46	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:PRO:O	1:A:285:ASP:HB2	2.08	0.53
1:B:65:PRO:HB3	1:B:93:LEU:HD13	1.93	0.50
1:A:42:LYS:NZ	4:A:503:HOH:O	2.43	0.49
1:A:258:THR:O	1:A:258:THR:HG22	2.12	0.47
1:B:281:LEU:HB3	1:B:288:ALA:O	2.15	0.47
1:B:204:HIS:CE1	1:B:206:ALA:HB2	2.51	0.46
1:A:73:LEU:HD22	1:A:93:LEU:HG	1.97	0.45
1:B:36:ILE:HD13	1:B:36:ILE:HG21	1.53	0.45
3:A:403:GOL:H32	4:A:547:HOH:O	2.15	0.43
1:A:277:TYR:HB2	1:A:294:MET:HG3	1.99	0.43
1:A:309:ARG:HB3	4:A:524:HOH:O	2.19	0.43
1:A:231:ALA:O	1:A:234:ASP:N	2.30	0.42
1:B:113:MET:HE1	1:B:212:ALA:HB3	2.01	0.42
1:B:65:PRO:HG3	1:B:89:PHE:HB2	2.00	0.41
1:A:229:GLU:CG	1:A:230:ALA:H	2.33	0.41
1:B:91:THR:HG23	1:B:92:ARG:N	2.35	0.41
1:A:249:ALA:C	1:A:250:ILE:HD12	2.40	0.41
1:B:112:THR:OG1	1:B:135[A]:HIS:HD2	1.98	0.41
1:B:118:GLY:O	1:B:133:GLY:N	2.54	0.41
1:B:60:THR:HB	1:B:86:PHE:CZ	2.56	0.40
1:A:250:ILE:CD1	1:A:250:ILE:N	2.84	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	273/296 (92%)	253 (93%)	16 (6%)	4 (2%)	12	9
1	B	272/296 (92%)	254 (93%)	14 (5%)	4 (2%)	12	9
All	All	545/592 (92%)	507 (93%)	30 (6%)	8 (2%)	12	9

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	118	GLY
1	A	230	ALA
1	B	67	PRO
1	B	267	GLU
1	A	117	GLY
1	A	228	SER
1	B	119	GLY
1	B	65	PRO

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	193/209 (92%)	183 (95%)	10 (5%)	27	32
1	B	192/209 (92%)	185 (96%)	7 (4%)	40	50
All	All	385/418 (92%)	368 (96%)	17 (4%)	34	40

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	ARG
1	A	51	ASN
1	A	190	ILE
1	A	222[A]	GLN
1	A	222[B]	GLN
1	A	235	VAL
1	A	243	ARG
1	A	264	ILE
1	A	272	LEU
1	A	309	ARG
1	B	29	VAL
1	B	67	PRO
1	B	195	GLN
1	B	257	ASP
1	B	261	LEU

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Mol	Chain	Res	Type
1	B	272	LEU
1	B	306	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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	$\# Z > 2$
3	GOL	A	402	-	5,5,5	0.30	0	5,5,5	0.73	0
3	GOL	A	403	-	5,5,5	0.54	0	5,5,5	1.27	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	A	402	-	-	0/4/4/4	0/0/0/0
3	GOL	A	403	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	403	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

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, 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	272/296 (91%)	0.13	14 (5%) 29 27	34, 47, 83, 118	0
1	B	271/296 (91%)	-0.07	10 (3%) 42 40	32, 46, 84, 106	0
All	All	543/592 (91%)	0.03	24 (4%) 35 33	32, 46, 84, 118	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	233	ALA	9.0
1	A	232	ALA	7.7
1	B	269	GLY	4.7
1	B	270	LEU	4.5
1	A	309	ARG	4.0
1	A	259	ARG	3.9
1	B	197	ARG	3.5
1	A	236	ALA	3.4
1	A	118	GLY	3.3
1	B	308	ALA	3.1
1	B	271	PRO	2.9
1	B	25	GLU	2.7
1	A	308	ALA	2.7
1	A	191	ALA	2.7
1	A	230	ALA	2.4
1	A	267	GLU	2.3
1	A	119	GLY	2.3
1	A	234	ASP	2.2
1	B	268	ALA	2.2
1	B	266	ARG	2.2
1	A	235	VAL	2.1
1	A	229	GLU	2.1
1	B	66	ARG	2.0
1	B	259	ARG	2.0

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(\AA^2)	Q<0.9
3	GOL	A	403	6/6	0.86	1.02	18.08	38,43,46,57	6
2	ZN	B	401	1/1	1.00	0.20	4.81	40,40,40,40	0
3	GOL	A	402	6/6	0.95	0.22	4.67	56,67,72,79	0
2	ZN	A	401	1/1	1.00	0.20	3.60	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.