



# wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 3WOO  
Title : Crystal structure of the DAP BII hexapeptide complex I  
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Deposited on : 2013-12-29  
Resolution : 1.80 Å(reported)

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We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

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<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>.

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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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

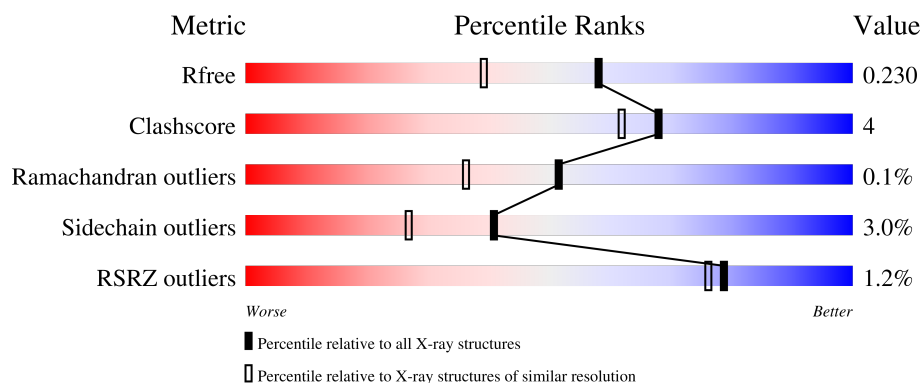
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
$R_{free}$	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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  $\geq 3$ , 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	698	<div> <div>92%</div> <div>7%</div> </div>
1	B	698	<div> <div>2%</div> <div>87%</div> <div>12%</div> </div>
2	C	6	<div> <div>17%</div> <div>33%</div> <div>50%</div> <div>17%</div> </div>
2	D	6	<div> <div>17%</div> <div>67%</div> <div>17%</div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11870 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 dipeptidyl aminopeptidase BII.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	697	Total	C	N	O	S	0	0	0
			5366	3395	935	1017	19			
1	B	697	Total	C	N	O	S	0	0	0
			5366	3395	935	1017	19			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	86	ALA	HIS	engineered mutation	UNP V5YM14
B	86	ALA	HIS	engineered mutation	UNP V5YM14

- Molecule 2 is a protein called Angiotensin II.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	5	Total	C	N	O	0	0	0
			44	31	7	6			
2	D	5	Total	C	N	O	0	0	0
			44	31	7	6			

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

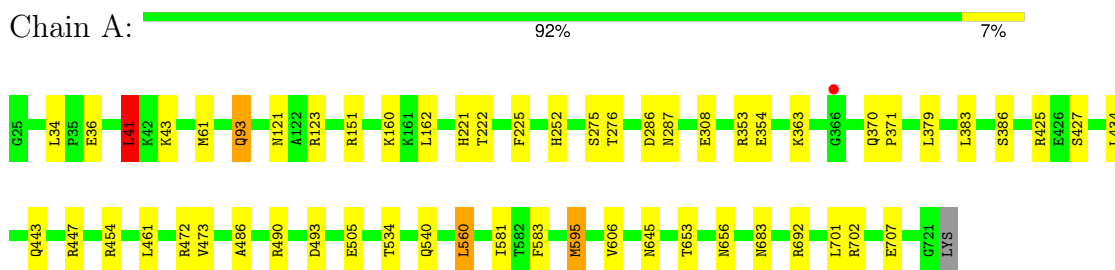
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	564	Total 564	O 564	0	0
5	B	404	Total 404	O 404	0	0
5	C	3	Total 3	O 3	0	0
5	D	3	Total 3	O 3	0	0

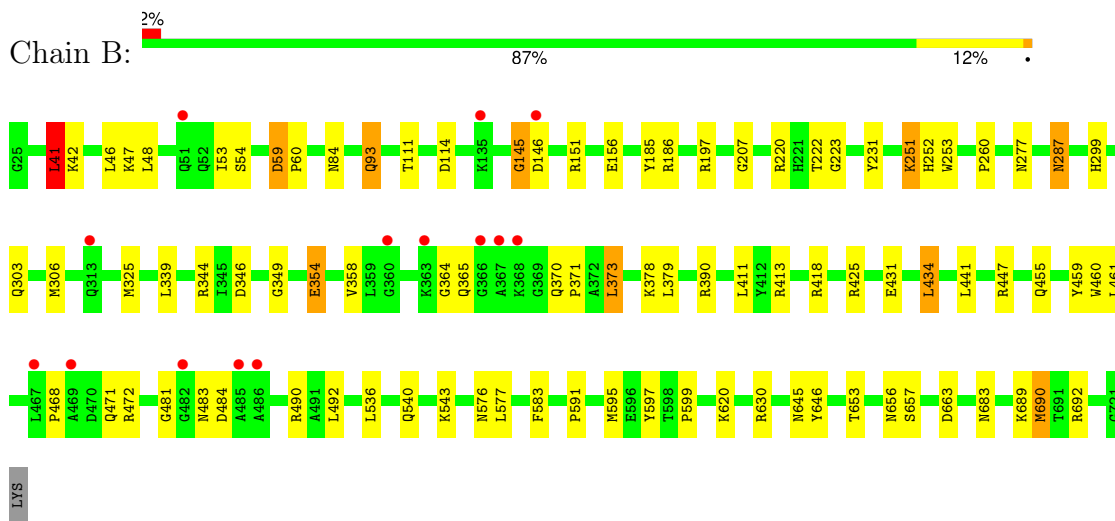
### 3 Residue-property plots

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.

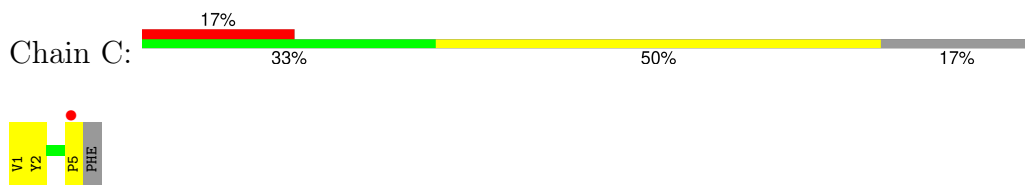
#### • Molecule 1: dipeptidyl aminopeptidase BII



#### • Molecule 1: dipeptidyl aminopeptidase BII



#### • Molecule 2: Angiotensin II



#### • Molecule 2: Angiotensin II





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	121.70Å 121.70Å 218.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.92 – 1.80 39.89 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (39.92-1.80) 99.5 (39.89-1.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.25 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.182 , 0.224 0.191 , 0.230	Depositor DCC
$R_{free}$ test set	7598 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.3	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 45.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	11870	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 23.37 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.7213e-03.*

<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, 0.333 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, 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	0.97	0/5483	0.99	14/7432 (0.2%)
1	B	0.90	1/5483 (0.0%)	0.95	14/7432 (0.2%)
2	C	1.44	0/46	1.19	0/63
2	D	1.53	0/46	1.24	0/63
All	All	0.94	1/11058 (0.0%)	0.97	28/14990 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	185	TYR	CE1-CZ	-5.41	1.31	1.38

The worst 5 of 28 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	325	MET	CG-SD-CE	-10.72	83.05	100.20
1	A	41	LEU	CA-CB-CG	7.62	132.84	115.30
1	B	692	ARG	NE-CZ-NH2	-7.54	116.53	120.30
1	B	690	MET	CG-SD-CE	-7.35	88.43	100.20
1	B	186	ARG	NE-CZ-NH1	7.04	123.82	120.30

There are no chirality outliers.

There are no planarity outliers.

### 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	5366	0	5282	30	0
1	B	5366	0	5282	46	0
2	C	44	0	45	4	0
2	D	44	0	45	0	0
3	A	48	0	64	7	0
3	B	24	0	32	3	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	A	564	0	0	13	1
5	B	404	0	0	3	0
5	C	3	0	0	1	0
5	D	3	0	0	0	0
All	All	11870	0	10750	82	1

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.

The worst 5 of 82 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:252:HIS:ND1	3:A:802:GOL:O1	1.89	1.04
3:A:805:GOL:H12	5:A:1026:HOH:O	1.67	0.94
1:A:252:HIS:HD1	3:A:802:GOL:HO1	1.17	0.90
1:A:707:GLU:OE2	5:A:1448:HOH:O	1.98	0.81
1:A:308:GLU:OE1	5:A:1452:HOH:O	2.06	0.73

All (1) 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)
5:A:1451:HOH:O	5:A:1457:HOH:O[4_555]	1.70	0.50

## 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	695/698 (100%)	685 (99%)	10 (1%)	0	100	100
1	B	695/698 (100%)	667 (96%)	26 (4%)	2 (0%)	41	27
2	C	3/6 (50%)	2 (67%)	1 (33%)	0	100	100
2	D	3/6 (50%)	2 (67%)	1 (33%)	0	100	100
All	All	1396/1408 (99%)	1356 (97%)	38 (3%)	2 (0%)	51	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	145	GLY
1	B	364	GLY

### 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	540/541 (100%)	527 (98%)	13 (2%)	49	36
1	B	540/541 (100%)	521 (96%)	19 (4%)	36	21
2	C	5/6 (83%)	5 (100%)	0	100	100
2	D	5/6 (83%)	4 (80%)	1 (20%)	1	0
All	All	1090/1094 (100%)	1057 (97%)	33 (3%)	41	27

5 of 33 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	461	LEU
1	B	492	LEU
2	D	2	TYR
1	A	560	LEU
1	A	540	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32

such sidechains are listed below:

Mol	Chain	Res	Type
1	B	656	ASN
1	B	665	HIS
1	A	645	ASN
1	A	585	ASN
2	C	4	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 4 are monoatomic - leaving 12 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	801	-	5,5,5	0.27	0	5,5,5	0.79	0
3	GOL	B	803	-	5,5,5	0.72	0	5,5,5	0.88	0
3	GOL	A	803	-	5,5,5	0.53	0	5,5,5	0.73	0
3	GOL	A	802	-	5,5,5	0.53	0	5,5,5	1.78	1 (20%)
3	GOL	A	806	-	5,5,5	0.33	0	5,5,5	0.38	0
3	GOL	B	802	-	5,5,5	0.35	0	5,5,5	0.48	0
3	GOL	A	804	-	5,5,5	0.63	0	5,5,5	0.95	0
3	GOL	B	801	-	5,5,5	0.66	0	5,5,5	1.18	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	GOL	B	804	-	5,5,5	0.48	0	5,5,5	0.44	0
3	GOL	A	805	-	5,5,5	2.33	1 (20%)	5,5,5	1.79	2 (40%)
3	GOL	A	808	-	5,5,5	0.46	0	5,5,5	1.09	1 (20%)
3	GOL	A	807	-	5,5,5	0.60	0	5,5,5	0.46	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	801	-	-	0/4/4/4	-
3	GOL	B	803	-	-	0/4/4/4	-
3	GOL	A	803	-	-	2/4/4/4	-
3	GOL	A	802	-	-	2/4/4/4	-
3	GOL	A	806	-	-	2/4/4/4	-
3	GOL	B	802	-	-	4/4/4/4	-
3	GOL	A	804	-	-	1/4/4/4	-
3	GOL	B	801	-	-	0/4/4/4	-
3	GOL	B	804	-	-	0/4/4/4	-
3	GOL	A	805	-	-	2/4/4/4	-
3	GOL	A	808	-	-	2/4/4/4	-
3	GOL	A	807	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	805	GOL	O2-C2	4.38	1.56	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	802	GOL	O1-C1-C2	-3.58	94.24	110.38
3	A	805	GOL	O2-C2-C3	2.70	120.37	109.18
3	A	805	GOL	O3-C3-C2	2.40	121.17	110.38
3	A	808	GOL	O3-C3-C2	-2.10	100.94	110.38

There are no chirality outliers.

5 of 17 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	802	GOL	O1-C1-C2-C3
3	A	803	GOL	O1-C1-C2-C3
3	A	805	GOL	C1-C2-C3-O3
3	A	808	GOL	C1-C2-C3-O3
3	A	808	GOL	O2-C2-C3-O3

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	803	GOL	3	0
3	A	802	GOL	3	0
3	A	806	GOL	2	0
3	A	805	GOL	2	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, 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	697/698 (99%)	-0.43	1 (0%) 95 93	10, 21, 35, 51	0
1	B	697/698 (99%)	-0.08	14 (2%) 65 61	13, 29, 50, 73	0
2	C	5/6 (83%)	1.40	1 (20%) 1 0	15, 19, 35, 55	0
2	D	5/6 (83%)	1.44	1 (20%) 1 0	17, 19, 35, 55	0
All	All	1404/1408 (99%)	-0.24	17 (1%) 79 76	10, 25, 46, 73	0

The worst 5 of 17 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	469	ALA	6.6
1	B	363	LYS	3.6
2	D	5	PRO	3.5
2	C	5	PRO	3.3
1	B	482	GLY	2.7

### 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 monosaccharides 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, 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	B-factors( $\text{\AA}^2$ )	Q<0.9
3	GOL	A	805	6/6	0.78	0.22	25,34,37,43	0
3	GOL	A	804	6/6	0.84	0.17	34,42,49,59	0
3	GOL	B	802	6/6	0.84	0.19	46,53,60,60	0
3	GOL	A	807	6/6	0.85	0.15	38,46,54,54	0
3	GOL	A	806	6/6	0.88	0.21	39,55,56,65	0
3	GOL	B	803	6/6	0.88	0.21	33,44,46,48	0
4	ZN	B	805	1/1	0.89	0.05	51,51,51,51	0
3	GOL	B	804	6/6	0.91	0.17	37,42,45,46	0
3	GOL	A	803	6/6	0.92	0.12	35,36,40,42	0
3	GOL	A	808	6/6	0.94	0.12	21,35,38,47	0
3	GOL	B	801	6/6	0.96	0.09	19,22,23,25	0
3	GOL	A	802	6/6	0.97	0.11	28,33,39,39	0
3	GOL	A	801	6/6	0.97	0.07	18,19,20,26	0
4	ZN	A	809	1/1	0.98	0.04	32,32,32,32	0
4	ZN	A	810	1/1	0.99	0.11	48,48,48,48	0
4	ZN	B	806	1/1	0.99	0.10	49,49,49,49	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.