



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 5, 2017 – 07:54 PM EST

PDB ID : 5MJM
Title : Single-shot pink beam serial crystallography: Phycocyanin (Five chips merged)
Authors : Meents, A.; Oberthuer, D.; Lieske, J.; Srajer, V.; Sarrou, I.
Deposited on : unknown
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 : 1.7.2 (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

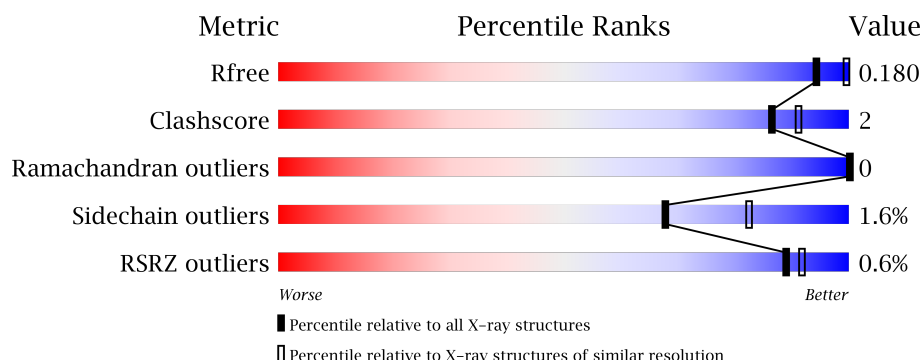
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

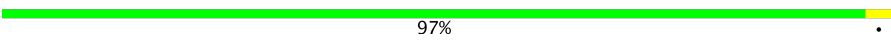
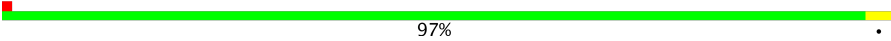
The reported resolution of this entry is 2.30 Å.

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}	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (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	162	 97%
2	B	172	 97%

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 2 is a protein called C-phycocyanin beta chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	172	Total	C	H	N	O	S	0	0	0
			2543	788	1270	228	249	8			

- # CYC

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 80	C 33	H 37	N 4	O 6	0	0
3	B	1	Total 80	C 33	H 37	N 4	O 6	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	H	N	O	0	0
			81	33	38	4	6		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	68	Total	O	0	0
			68	68		
4	B	59	Total	O	0	0
			59	59		

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.

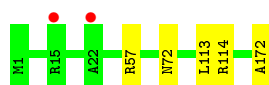
- Molecule 1: C-phycocyanin alpha chain

Chain A:  97% .



- Molecule 2: C-phycocyanin beta chain

Chain B:  97% .



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	187.80Å 187.80Å 60.70Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	14.92 – 2.30 27.42 – 2.30	Depositor EDS
% Data completeness (in resolution range)	66.8 (14.92-2.30) 66.9 (27.42-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.38 (at 2.31Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.139 , 0.181 0.130 , 0.180	Depositor DCC
R_{free} test set	1223 reflections (10.06%)	DCC
Wilson B-factor (Å ²)	12.0	Xtriage
Anisotropy	0.287	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 35.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5339	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.79% 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: CYC, MEN

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	0.34	0/1252	0.44	0/1698
2	B	0.32	0/1277	0.44	0/1727
All	All	0.33	0/2529	0.44	0/3425

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	1225	1203	1197	1	0
2	B	1273	1270	1271	4	0
3	A	43	37	37	4	0
3	B	86	75	73	7	0
4	A	68	0	0	0	0
4	B	59	0	0	1	0
All	All	2754	2585	2578	13	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 2.

All (13) 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)
3:B:201:CYC:HC	3:B:201:CYC:HMD2	1.59	0.67
2:B:57:ARG:NH1	4:B:305:HOH:O	2.33	0.61
3:B:202:CYC:NC	3:B:202:CYC:HMD2	2.21	0.55
3:B:201:CYC:NC	3:B:201:CYC:HMD2	2.22	0.54
2:B:114:ARG:NH2	2:B:172:ALA:O	2.33	0.54
3:B:201:CYC:NB	3:B:201:CYC:HMA1	2.22	0.54
3:A:501:CYC:NB	3:A:501:CYC:HMA1	2.26	0.50
2:B:72:MEN:HE22	3:B:201:CYC:HBD2	1.92	0.50
3:A:501:CYC:HBC2	3:A:501:CYC:HMC1	1.94	0.49
2:B:113:LEU:HD13	3:B:201:CYC:HMB3	1.96	0.47
3:B:202:CYC:HMA1	3:B:202:CYC:NB	2.30	0.47
1:A:84:CYS:HA	3:A:501:CYC:HHD	2.01	0.42
3:A:501:CYC:HMD2	3:A:501:CYC:NC	2.35	0.42

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	161/162 (99%)	158 (98%)	3 (2%)	0	100	100
2	B	169/172 (98%)	167 (99%)	2 (1%)	0	100	100
All	All	330/334 (99%)	325 (98%)	5 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	123/123 (100%)	119 (97%)	4 (3%)	43	59
2	B	127/127 (100%)	127 (100%)	0	100	100
All	All	250/250 (100%)	246 (98%)	4 (2%)	68	82

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	30	ARG
1	A	38	MET
1	A	140	HIS

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 ⓘ

1 non-standard protein/DNA/RNA residue is 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	MEN	B	72	2	8,8,9	1.21	1 (12%)	8,9,11	0.94	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	MEN	B	72	2	-	0/6/8/10	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	72	MEN	CA-C	2.60	1.53	1.50

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 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	72	MEN	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
3	CYC	A	501	1	38,46,46	3.57	18 (47%)	46,67,67	1.78	9 (19%)
3	CYC	B	201	2	38,46,46	3.62	18 (47%)	46,67,67	1.94	13 (28%)
3	CYC	B	202	2	38,46,46	3.62	19 (50%)	46,67,67	1.72	12 (26%)

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	CYC	A	501	1	-	3/21/74/74	0/4/4/4
3	CYC	B	201	2	-	3/21/74/74	0/4/4/4
3	CYC	B	202	2	-	3/21/74/74	0/4/4/4

All (55) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	CYC	CHA-C1A	-5.27	1.30	1.35
3	B	201	CYC	C2C-C1C	-4.41	1.47	1.52
3	B	201	CYC	CHA-C1A	-4.33	1.31	1.35
3	B	202	CYC	CHA-C1A	-4.10	1.31	1.35
3	B	202	CYC	C2C-C1C	-3.73	1.48	1.52
3	A	501	CYC	C2C-C1C	-3.17	1.49	1.52
3	A	501	CYC	CHB-C1B	-2.69	1.31	1.37
3	B	201	CYC	CHB-C1B	-2.56	1.31	1.37
3	A	501	CYC	OC-C1C	-2.43	1.18	1.23
3	A	501	CYC	OB-C4B	-2.42	1.19	1.23
3	B	202	CYC	OB-C4B	-2.35	1.19	1.23
3	B	202	CYC	CHB-C1B	-2.32	1.32	1.37
3	B	201	CYC	OB-C4B	-2.29	1.19	1.23
3	B	201	CYC	OC-C1C	-2.20	1.19	1.23
3	B	202	CYC	OC-C1C	-2.16	1.19	1.23
3	B	202	CYC	CHD-C4C	-2.12	1.32	1.38
3	B	201	CYC	C3B-C2B	2.28	1.41	1.36
3	B	201	CYC	C1B-C2B	2.32	1.49	1.45
3	A	501	CYC	C4D-CHA	2.36	1.52	1.45
3	A	501	CYC	C1B-C2B	2.36	1.49	1.45
3	A	501	CYC	C3B-C2B	2.45	1.42	1.36
3	B	202	CYC	C1B-C2B	2.48	1.49	1.45
3	B	201	CYC	C4D-CHA	2.50	1.52	1.45
3	B	202	CYC	C4D-CHA	2.69	1.53	1.45
3	B	202	CYC	C1D-CHD	2.82	1.53	1.45
3	B	202	CYC	C3B-C2B	2.88	1.42	1.36
3	A	501	CYC	C1D-CHD	2.97	1.54	1.45
3	B	201	CYC	C1D-CHD	3.06	1.54	1.45
3	A	501	CYC	C1A-C2A	3.80	1.52	1.45
3	B	201	CYC	C1A-C2A	4.16	1.52	1.45
3	B	202	CYC	C1A-C2A	4.36	1.52	1.45
3	A	501	CYC	C2A-C3A	4.48	1.46	1.36
3	B	202	CYC	C2A-C3A	4.61	1.46	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	201	CYC	C2A-C3A	4.85	1.47	1.36
3	A	501	CYC	C4A-NA	5.10	1.48	1.36
3	B	202	CYC	C4A-NA	5.31	1.48	1.36
3	B	202	CYC	CHB-C4A	5.35	1.53	1.40
3	A	501	CYC	CHB-C4A	5.38	1.53	1.40
3	B	201	CYC	CHB-C4A	5.41	1.53	1.40
3	B	201	CYC	C4A-NA	5.41	1.49	1.36
3	A	501	CYC	C1A-NA	5.46	1.50	1.38
3	B	202	CYC	C1A-NA	5.65	1.50	1.38
3	B	201	CYC	C1A-NA	5.87	1.51	1.38
3	A	501	CYC	C4C-NC	6.23	1.50	1.37
3	B	202	CYC	C4C-NC	6.51	1.51	1.37
3	B	201	CYC	C4C-NC	6.65	1.51	1.37
3	A	501	CYC	C4B-NB	7.30	1.53	1.37
3	B	201	CYC	C4B-NB	7.48	1.53	1.37
3	B	202	CYC	C4B-NB	7.57	1.53	1.37
3	A	501	CYC	C1B-NB	8.77	1.53	1.37
3	B	201	CYC	C1B-NB	8.80	1.53	1.37
3	B	202	CYC	C1B-NB	9.06	1.53	1.37
3	B	201	CYC	C1C-NC	9.63	1.49	1.37
3	B	202	CYC	C1C-NC	9.66	1.49	1.37
3	A	501	CYC	C1C-NC	9.74	1.50	1.37

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	CYC	C3C-C4C-NC	-6.10	102.21	107.97
3	B	202	CYC	CAA-CBA-CGA	-2.99	107.55	112.66
3	A	501	CYC	CBD-CAD-C3D	-2.97	106.80	112.48
3	B	201	CYC	C3C-C4C-NC	-2.89	105.24	107.97
3	B	202	CYC	C2A-C1A-NA	-2.87	105.55	109.93
3	A	501	CYC	C1B-NB-C4B	-2.78	107.00	110.70
3	B	201	CYC	CAB-C3B-C2B	-2.61	122.96	127.49
3	B	202	CYC	CAC-C3C-C2C	-2.56	107.81	114.24
3	B	201	CYC	C2A-C1A-NA	-2.52	106.08	109.93
3	A	501	CYC	C1B-CHB-C4A	-2.44	121.93	128.08
3	B	201	CYC	CAC-C3C-C2C	-2.34	108.36	114.24
3	B	201	CYC	CBD-CAD-C3D	-2.34	108.01	112.48
3	B	202	CYC	C3C-C4C-NC	-2.33	105.77	107.97
3	B	202	CYC	C1B-NB-C4B	-2.32	107.62	110.70
3	B	201	CYC	C1B-NB-C4B	-2.30	107.64	110.70
3	B	201	CYC	CMC-C2C-C1C	-2.22	107.78	112.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	201	CYC	C1B-CHB-C4A	-2.08	122.85	128.08
3	B	202	CYC	CBD-CAD-C3D	-2.07	108.53	112.48
3	B	201	CYC	C1A-C2A-C3A	2.07	109.13	106.81
3	A	501	CYC	C1A-C2A-C3A	2.11	109.17	106.81
3	A	501	CYC	CMA-C3A-C4A	2.60	129.09	125.04
3	B	202	CYC	C1A-C2A-C3A	2.61	109.73	106.81
3	B	202	CYC	CMA-C3A-C4A	2.64	129.15	125.04
3	A	501	CYC	C1B-C2B-C3B	2.76	110.85	107.83
3	B	201	CYC	C1B-C2B-C3B	3.01	111.13	107.83
3	B	202	CYC	C1B-C2B-C3B	3.02	111.13	107.83
3	A	501	CYC	C4B-C3B-C2B	3.20	109.85	108.01
3	B	202	CYC	C4B-C3B-C2B	3.52	110.03	108.01
3	B	202	CYC	C2C-C3C-C4C	3.76	106.97	101.34
3	B	202	CYC	CAB-C3B-C4B	3.86	124.78	121.53
3	B	201	CYC	C4B-C3B-C2B	4.23	110.44	108.01
3	B	201	CYC	C2C-C3C-C4C	4.67	108.34	101.34
3	A	501	CYC	CAB-C3B-C4B	4.96	125.71	121.53
3	B	201	CYC	CAB-C3B-C4B	6.01	126.59	121.53

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	201	CYC	C1B-CHB-C4A-C3A
3	B	202	CYC	C1B-CHB-C4A-C3A
3	A	501	CYC	C1B-CHB-C4A-C3A
3	B	201	CYC	C1B-CHB-C4A-NA
3	A	501	CYC	C1B-CHB-C4A-NA
3	A	501	CYC	C4C-CHD-C1D-ND
3	B	202	CYC	C4C-CHD-C1D-ND
3	B	202	CYC	C1B-CHB-C4A-NA
3	B	201	CYC	C4C-CHD-C1D-ND

There are no ring outliers.

3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	CYC	4	0
3	B	201	CYC	5	0
3	B	202	CYC	2	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	162/162 (100%)	-0.59	0 100 100	11, 17, 33, 75	0
2	B	171/172 (99%)	-0.40	2 (1%) 79 82	13, 23, 39, 57	0
All	All	333/334 (99%)	-0.49	2 (0%) 89 92	11, 20, 38, 75	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	22	ALA	2.1
2	B	15	ARG	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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	MEN	B	72	9/10	0.97	0.10	-	15,26,36,36	0

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	CYC	B	201	43/43	0.92	0.15	0.85	13,26,41,43	0
3	CYC	A	501	43/43	0.97	0.10	0.10	2,11,20,22	0
3	CYC	B	202	43/43	0.94	0.13	0.06	15,23,30,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.