



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 04:43 AM GMT

PDB ID : 2NZJ
Title : The crystal structure of REM1 in complex with GDP
Authors : Turnbull, A.P.; Papagrigoriou, E.; Ugochukwu, E.; Elkins, J.M.; Soundararajan, M.; Yang, X.; Gorrec, F.; Umeano, C.; Salah, E.; Burgess, N.; Johansson, C.; Berridge, G.; Gileadi, O.; Bray, J.; Marsden, B.; Watts, S.; von Delft, F.; Weigelt, J.; Edwards, A.; Arrowsmith, C.H.; Sundstrom, M.; Doyle, D.; Structural Genomics Consortium (SGC)
Deposited on : 2006-11-23
Resolution : 2.50 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

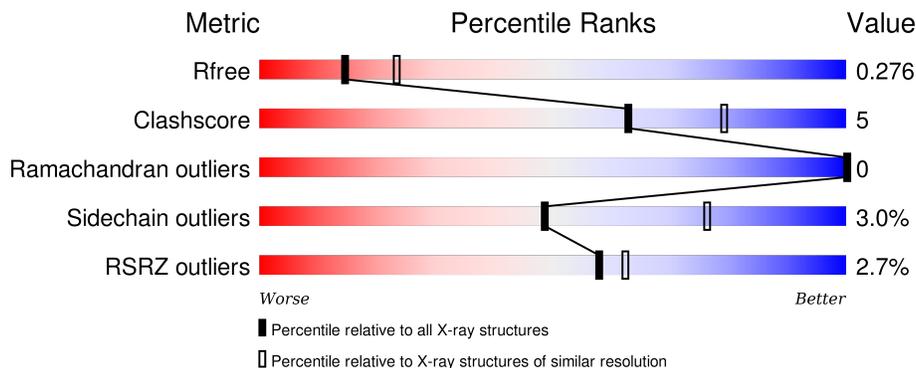
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	175	<p>2% 77% 11% • 11%</p>
1	B	175	<p>3% 77% 14% 9%</p>
1	C	175	<p>3% 81% 9% • 9%</p>
1	D	175	<p>2% 73% 14% • 11%</p>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4785 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 GTP-binding protein REM 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	155	1140	722	196	217	5	0	0	0
1	B	160	1180	744	210	221	5	0	0	0
1	C	159	1185	747	207	226	5	0	0	0
1	D	155	1146	721	200	220	5	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	CLONING ARTIFACT	UNP O75628
A	0	MET	-	CLONING ARTIFACT	UNP O75628
B	-1	SER	-	CLONING ARTIFACT	UNP O75628
B	0	MET	-	CLONING ARTIFACT	UNP O75628
C	-1	SER	-	CLONING ARTIFACT	UNP O75628
C	0	MET	-	CLONING ARTIFACT	UNP O75628
D	-1	SER	-	CLONING ARTIFACT	UNP O75628
D	0	MET	-	CLONING ARTIFACT	UNP O75628

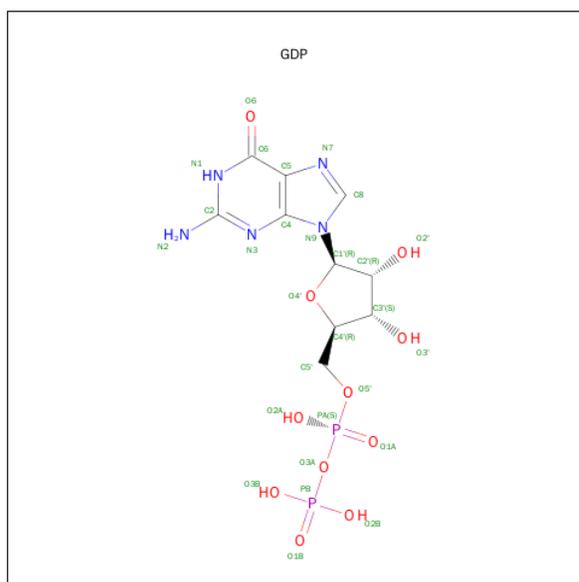
- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total 1	Mg 1	0	0
2	A	1	Total 1	Mg 1	0	0
2	D	1	Total 1	Mg 1	0	0
2	C	1	Total 1	Mg 1	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 4 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O P 28 10 5 11 2	0	0
4	B	1	Total C N O P 28 10 5 11 2	0	0
4	C	1	Total C N O P 28 10 5 11 2	0	0
4	D	1	Total C N O P 28 10 5 11 2	0	0

- Molecule 5 is water.

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

Continued on next page...

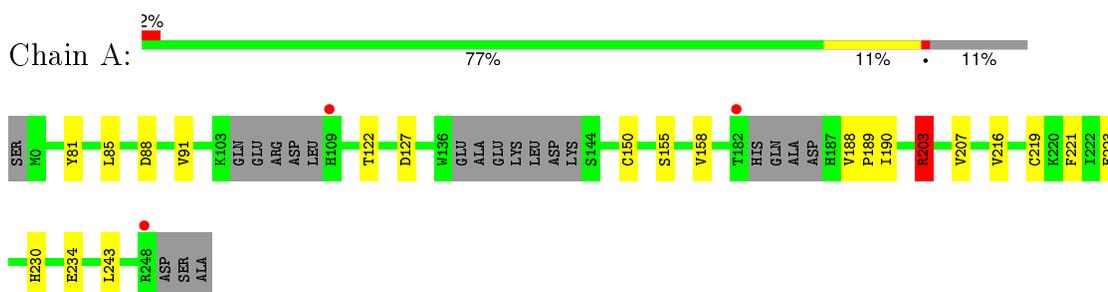
Continued from previous page...

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	4	Total O 4 4	0	0
5	D	4	Total O 4 4	0	0

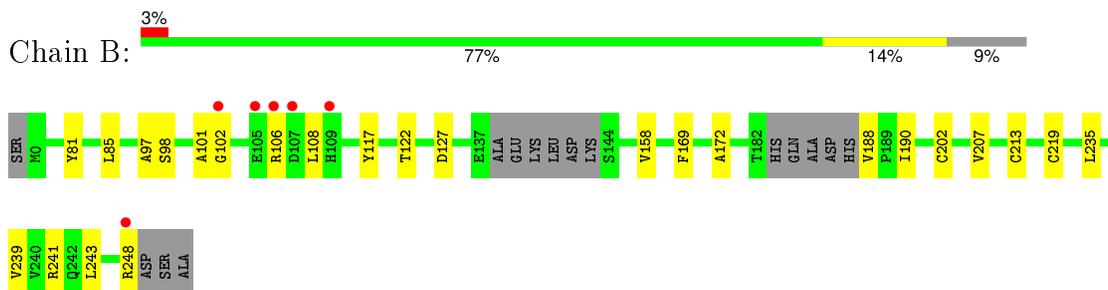
3 Residue-property plots [i](#)

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 ($RSRZ > 2$). 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.

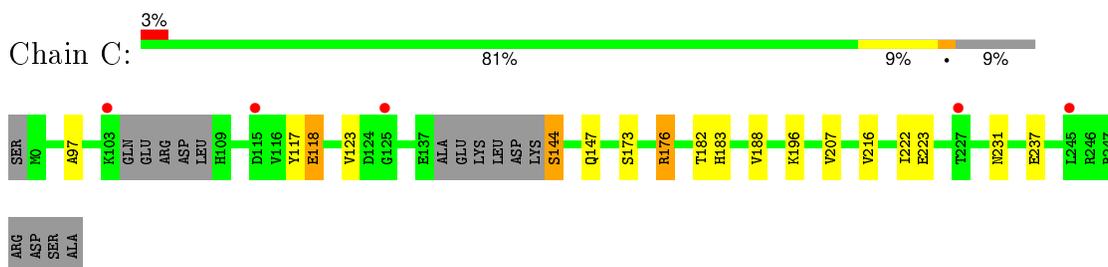
- Molecule 1: GTP-binding protein REM 1



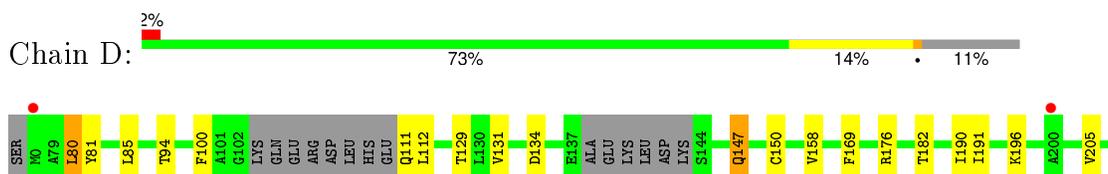
- Molecule 1: GTP-binding protein REM 1



- Molecule 1: GTP-binding protein REM 1



- Molecule 1: GTP-binding protein REM 1



E209	C213	K220	L235	F236	V239	L243	R244	L245	R246	ARG	ASP	SER	ALA
------	------	------	------	------	------	------	------	------	------	-----	-----	-----	-----

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	43.64Å 102.28Å 165.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.00 – 2.50 30.03 – 2.51	Depositor EDS
% Data completeness (in resolution range)	97.7 (29.00-2.50) 97.7 (30.03-2.51)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.58 (at 2.51Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.225 , 0.280 0.224 , 0.276	Depositor DCC
R_{free} test set	1284 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	47.6	Xtrriage
Anisotropy	0.347	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 24.7	EDS
Estimated twinning fraction	No twinning to report.	Xtrriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Outliers	2 of 25472 reflections (0.008%)	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4785	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 60.23 % 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 1.5684e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: GDP, MG, CL

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	0.69	1/1152 (0.1%)	0.71	1/1563 (0.1%)
1	B	0.64	1/1193 (0.1%)	0.70	1/1617 (0.1%)
1	C	0.60	0/1200	0.67	0/1627
1	D	0.57	0/1160	0.67	0/1573
All	All	0.63	2/4705 (0.0%)	0.69	2/6380 (0.0%)

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
1	D	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	202	CYS	CB-SG	-5.69	1.72	1.81
1	A	234	GLU	CG-CD	5.26	1.59	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	203	ARG	NE-CZ-NH1	-6.56	117.02	120.30
1	B	241	ARG	NE-CZ-NH1	5.91	123.26	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	101	ALA	Peptide
1	D	112	LEU	Peptide

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	1140	0	1102	12	0
1	B	1180	0	1148	13	0
1	C	1185	0	1147	11	0
1	D	1146	0	1098	17	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	28	0	12	0	0
4	B	28	0	12	0	0
4	C	28	0	12	1	0
4	D	28	0	12	1	0
5	A	4	0	0	0	0
5	B	4	0	0	0	0
5	C	4	0	0	0	0
5	D	4	0	0	0	0
All	All	4785	0	4543	49	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 5.

The worst 5 of 49 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:D:80:LEU:HD13	1:D:129:THR:HG22	1.69	0.73
1:D:176:ARG:NH2	1:D:213:CYS:O	2.31	0.64
1:D:191:ILE:N	1:D:191:ILE:HD13	2.16	0.61
1:D:147:GLN:HA	1:D:147:GLN:HE21	1.66	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:123:VAL:HG22	1:C:237:GLU:HG2	1.83	0.60

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	147/175 (84%)	142 (97%)	5 (3%)	0	100	100
1	B	154/175 (88%)	149 (97%)	5 (3%)	0	100	100
1	C	153/175 (87%)	152 (99%)	1 (1%)	0	100	100
1	D	149/175 (85%)	147 (99%)	2 (1%)	0	100	100
All	All	603/700 (86%)	590 (98%)	13 (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	115/148 (78%)	112 (97%)	3 (3%)	54	81
1	B	117/148 (79%)	116 (99%)	1 (1%)	84	95
1	C	120/148 (81%)	115 (96%)	5 (4%)	36	62
1	D	114/148 (77%)	109 (96%)	5 (4%)	35	60

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	466/592 (79%)	452 (97%)	14 (3%)	48 76

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	176	ARG
1	C	182	THR
1	D	150	CYS
1	C	144	SER
1	D	147	GLN

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

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 6 are monoatomic - leaving 4 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
4	GDP	A	1001	2	23,30,30	1.07	2 (8%)	30,47,47	1.97	7 (23%)
4	GDP	B	1002	2	23,30,30	1.10	2 (8%)	30,47,47	2.00	10 (33%)
4	GDP	C	1003	2	23,30,30	1.05	2 (8%)	30,47,47	1.87	8 (26%)
4	GDP	D	1004	2	23,30,30	1.13	3 (13%)	30,47,47	1.95	8 (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
4	GDP	A	1001	2	-	0/12/32/32	0/3/3/3
4	GDP	B	1002	2	-	0/12/32/32	0/3/3/3
4	GDP	C	1003	2	-	0/12/32/32	0/3/3/3
4	GDP	D	1004	2	-	0/12/32/32	0/3/3/3

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1004	GDP	O4'-C1'	2.15	1.43	1.41
4	D	1004	GDP	C5-C4	2.63	1.46	1.40
4	A	1001	GDP	C5-C4	2.64	1.46	1.40
4	C	1003	GDP	C5-C4	2.66	1.46	1.40
4	A	1001	GDP	C6-C5	2.69	1.46	1.41

The worst 5 of 33 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1004	GDP	C5-C6-N1	-5.28	116.37	123.59
4	C	1003	GDP	C5-C6-N1	-4.35	117.63	123.59
4	A	1001	GDP	C5-C6-N1	-4.28	117.74	123.59
4	B	1002	GDP	C5-C6-N1	-4.25	117.78	123.59
4	B	1002	GDP	PA-O3A-PB	-4.13	118.83	132.67

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
4	C	1003	GDP	1	0
4	D	1004	GDP	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	155/175 (88%)	-0.14	3 (1%) 70 73	10, 25, 35, 41	0
1	B	160/175 (91%)	0.04	6 (3%) 44 49	10, 27, 36, 42	0
1	C	159/175 (90%)	0.04	5 (3%) 52 57	5, 23, 36, 41	0
1	D	155/175 (88%)	0.01	3 (1%) 70 73	4, 24, 35, 39	0
All	All	629/700 (89%)	-0.01	17 (2%) 58 62	4, 25, 36, 42	0

The worst 5 of 17 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	105	GLU	6.8
1	C	245	LEU	5.1
1	C	115	ASP	4.0
1	B	102	GLY	3.8
1	B	109	HIS	3.7

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 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
4	GDP	C	1003	28/28	0.97	0.08	-1.12	39,41,43,44	0
4	GDP	B	1002	28/28	0.97	0.09	-1.38	34,37,38,39	0
4	GDP	A	1001	28/28	0.98	0.09	-1.44	29,36,37,41	0
4	GDP	D	1004	28/28	0.97	0.08	-1.50	38,41,45,46	0
3	CL	B	253	1/1	0.97	0.06	-1.76	45,45,45,45	0
3	CL	A	253	1/1	0.98	0.07	-2.36	50,50,50,50	0
2	MG	A	252	1/1	0.95	0.07	-	37,37,37,37	0
2	MG	C	252	1/1	0.97	0.04	-	49,49,49,49	0
2	MG	B	252	1/1	0.95	0.04	-	36,36,36,36	0
2	MG	D	252	1/1	0.94	0.03	-	50,50,50,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.