



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 02:33 PM GMT

PDB ID : 2CXX  
Title : Crystal structure of a probable GTP-binding protein engB  
Authors : Kukimoto-Niino, M.; Murayama, K.; Shirouzu, M.; Yokoyama, S.; RIKEN  
Structural Genomics/Proteomics Initiative (RSGI)  
Deposited on : 2005-07-01  
Resolution : 1.70 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

---

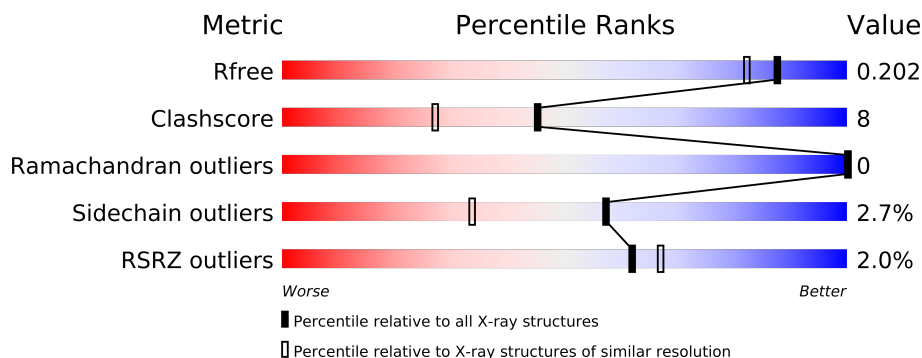
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.70 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))
$R_{free}$	66092	2456 (1.70-1.70)
Clashscore	79885	2929 (1.70-1.70)
Ramachandran outliers	78287	2878 (1.70-1.70)
Sidechain outliers	78261	2878 (1.70-1.70)
RSRZ outliers	66119	2456 (1.70-1.70)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	190	
1	B	190	
1	C	190	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5054 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Probable GTP-binding protein engB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	184	Total	C	N	O	Se	0	0	0
			1503	973	264	263	3			
1	B	176	Total	C	N	O	Se	0	0	0
			1443	937	249	254	3			
1	C	179	Total	C	N	O	Se	0	0	0
			1461	948	252	258	3			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP O57939
A	51	MSE	MET	MODIFIED RESIDUE	UNP O57939
A	57	MSE	MET	MODIFIED RESIDUE	UNP O57939
A	58	MSE	MET	MODIFIED RESIDUE	UNP O57939
B	1	MSE	MET	MODIFIED RESIDUE	UNP O57939
B	51	MSE	MET	MODIFIED RESIDUE	UNP O57939
B	57	MSE	MET	MODIFIED RESIDUE	UNP O57939
B	58	MSE	MET	MODIFIED RESIDUE	UNP O57939
C	1	MSE	MET	MODIFIED RESIDUE	UNP O57939
C	51	MSE	MET	MODIFIED RESIDUE	UNP O57939
C	57	MSE	MET	MODIFIED RESIDUE	UNP O57939
C	58	MSE	MET	MODIFIED RESIDUE	UNP O57939

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>11</sub>P<sub>2</sub>).



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

- Molecule 3 is water.

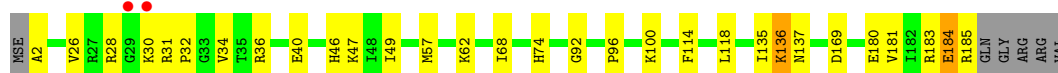
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	201	Total	O	0	0
			201	201		
3	B	198	Total	O	0	0
			198	198		
3	C	164	Total	O	0	0
			164	164		

### 3 Residue-property plots

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.

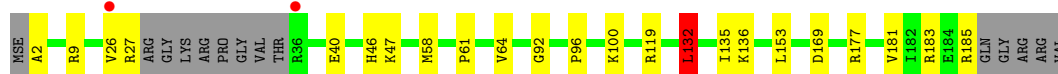
- Molecule 1: Probable GTP-binding protein engB

Chain A: 



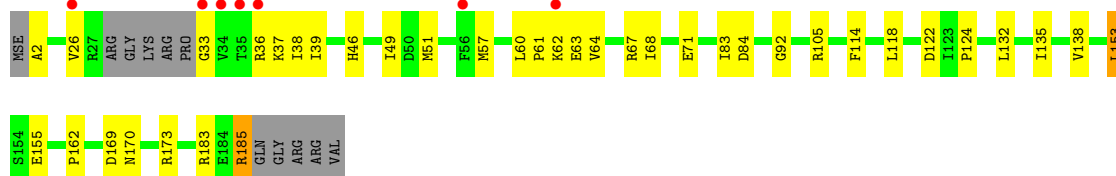
- Molecule 1: Probable GTP-binding protein engB

Chain B: 



- Molecule 1: Probable GTP-binding protein engB

Chain C: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	75.36Å 75.36Å 79.11Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.06 – 1.70 19.06 – 1.70	Depositor EDS
% Data completeness (in resolution range)	94.8 (19.06-1.70) 94.9 (19.06-1.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	8.16 (at 1.70Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.164 , 0.202 0.163 , 0.202	Depositor DCC
$R_{free}$ test set	5308 reflections (10.12%)	DCC
Wilson B-factor (Å <sup>2</sup> )	13.3	Xtriage
Anisotropy	0.308	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 36.2	EDS
Estimated twinning fraction	0.021 for -h,-k,l 0.044 for h,-h-k,-l 0.029 for -k,-h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 52458 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5054	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

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

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	0.84	0/1526	0.90	0/2045
1	B	0.88	0/1464	0.93	4/1961 (0.2%)
1	C	0.88	0/1482	0.92	2/1986 (0.1%)
All	All	0.87	0/4472	0.92	6/5992 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	9	ARG	NE-CZ-NH2	-6.22	117.19	120.30
1	C	169	ASP	CB-CG-OD1	6.19	123.87	118.30
1	B	132	LEU	CA-CB-CG	5.92	128.91	115.30
1	C	183	ARG	NE-CZ-NH1	5.80	123.20	120.30
1	B	177	ARG	NE-CZ-NH2	-5.67	117.47	120.30
1	B	177	ARG	NE-CZ-NH1	5.04	122.82	120.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1503	0	1583	21	0
1	B	1443	0	1514	14	0
1	C	1461	0	1533	42	0
2	A	28	0	12	0	0
2	B	28	0	12	0	0
2	C	28	0	12	0	0
3	A	201	0	0	6	0
3	B	198	0	0	2	0
3	C	164	0	0	5	0
All	All	5054	0	4666	77	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (77) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:26:VAL:HG22	3:A:1471:HOH:O	1.64	0.95
1:C:49:ILE:HG22	1:C:51:MSE:CE	2.12	0.79
1:A:181:VAL:O	1:A:184:GLU:HG3	1.94	0.66
1:B:40:GLU:CD	1:B:47:LYS:HD3	2.16	0.66
1:C:49:ILE:HG22	1:C:51:MSE:HE3	1.76	0.66
1:A:36:ARG:HG2	1:A:68:ILE:HD11	1.78	0.64
1:A:181:VAL:HA	1:A:184:GLU:HG2	1.80	0.64
1:C:61:PRO:HB2	1:C:64:VAL:HG23	1.80	0.63
1:C:122:ASP:O	1:C:185:ARG:NH2	2.31	0.63
1:C:170:ASN:HD21	1:C:173:ARG:HD3	1.64	0.62
1:C:57:MSE:HE3	1:C:60:LEU:HG	1.80	0.62
1:C:38:ILE:HG23	1:C:51:MSE:CE	2.30	0.62
1:A:74:HIS:ND1	3:A:1492:HOH:O	2.30	0.61
1:A:2:ALA:N	1:A:46:HIS:HD1	1.99	0.61
1:C:36:ARG:NH2	1:C:60:LEU:HD23	2.17	0.60
1:C:36:ARG:CZ	1:C:60:LEU:HD23	2.33	0.59
1:A:183:ARG:HD3	3:A:1455:HOH:O	2.02	0.59
1:C:92:GLY:HA3	1:C:135:ILE:HD11	1.85	0.59
1:C:170:ASN:ND2	1:C:173:ARG:HD3	2.19	0.58
1:B:2:ALA:N	1:B:46:HIS:HD1	2.02	0.58
1:A:136:LYS:HG3	1:A:137:ASN:N	2.19	0.57
1:B:40:GLU:OE2	1:B:47:LYS:HD3	2.04	0.57
1:C:38:ILE:HG23	1:C:51:MSE:HE1	1.86	0.57
1:C:2:ALA:N	1:C:46:HIS:HD1	2.04	0.56
1:B:92:GLY:HA3	1:B:135:ILE:HD11	1.87	0.56
1:B:136:LYS:HE2	3:B:2594:HOH:O	2.06	0.56
1:C:173:ARG:HG3	3:C:3464:HOH:O	2.07	0.55

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:181:VAL:O	1:B:185:ARG:HG2	2.07	0.54
1:C:26:VAL:CG1	1:C:39:ILE:HG21	2.38	0.54
1:B:26:VAL:O	1:B:27:ARG:HB2	2.06	0.54
1:C:170:ASN:HD21	1:C:173:ARG:HH11	1.56	0.54
1:C:124:PRO:HD3	1:C:185:ARG:NH2	2.24	0.53
1:A:181:VAL:HA	1:A:184:GLU:CG	2.39	0.53
1:B:26:VAL:O	1:B:27:ARG:CB	2.57	0.52
1:C:67:ARG:O	1:C:71:GLU:HG3	2.09	0.52
1:C:185:ARG:NH2	3:C:3518:HOH:O	2.42	0.52
1:C:33:GLY:O	1:C:37:LYS:HG2	2.10	0.52
1:C:36:ARG:NE	1:C:60:LEU:HD23	2.25	0.51
1:A:30:LYS:O	1:A:32:PRO:HD3	2.10	0.51
1:C:84:ASP:HA	1:C:185:ARG:HD2	1.94	0.50
1:C:170:ASN:HD21	1:C:173:ARG:NH1	2.10	0.50
1:C:153:LEU:O	1:C:153:LEU:HD22	2.12	0.50
1:C:105:ARG:CG	3:C:3490:HOH:O	2.60	0.50
1:C:83:ILE:O	1:C:185:ARG:NH1	2.45	0.49
1:A:180:GLU:O	1:A:184:GLU:HG2	2.13	0.49
1:A:96:PRO:O	1:A:100:LYS:HG3	2.12	0.49
1:B:96:PRO:O	1:B:100:LYS:HG3	2.12	0.49
1:C:49:ILE:HG22	1:C:51:MSE:HE1	1.94	0.49
1:C:114:PHE:CE1	1:C:118:LEU:HD11	2.47	0.49
1:C:36:ARG:HB3	1:C:68:ILE:HD11	1.94	0.48
1:B:132:LEU:HA	1:B:135:ILE:HD12	1.96	0.48
1:A:183:ARG:HG3	3:A:1559:HOH:O	2.14	0.47
1:C:38:ILE:HG23	1:C:51:MSE:HE2	1.96	0.46
1:C:132:LEU:HA	1:C:135:ILE:HD12	1.98	0.46
1:A:36:ARG:NH2	3:A:1601:HOH:O	2.49	0.46
1:C:38:ILE:CG2	1:C:51:MSE:HE1	2.45	0.45
1:C:62:LYS:HB3	1:C:62:LYS:HE2	1.69	0.44
1:C:26:VAL:HG12	1:C:39:ILE:HD13	2.00	0.44
1:C:124:PRO:HD3	1:C:185:ARG:CZ	2.48	0.44
1:A:57:MSE:HE3	3:A:1488:HOH:O	2.18	0.44
1:C:61:PRO:HB2	1:C:64:VAL:CG2	2.47	0.44
1:B:153:LEU:HD12	1:B:153:LEU:HA	1.89	0.43
1:A:40:GLU:HG2	1:A:49:ILE:HG12	2.00	0.43
1:C:60:LEU:HD13	1:C:61:PRO:O	2.18	0.43
1:A:114:PHE:CE2	1:A:118:LEU:HD11	2.53	0.43
1:C:138:VAL:HG21	1:C:162:PRO:HB3	2.00	0.43
1:A:92:GLY:HA3	1:A:135:ILE:HD11	2.01	0.42
1:A:40:GLU:HB3	1:A:47:LYS:HD2	2.01	0.42
1:B:119:ARG:NH2	3:B:2528:HOH:O	2.52	0.41

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:28:ARG:HD3	1:A:28:ARG:HA	1.84	0.41
1:C:155:GLU:HA	1:C:155:GLU:OE2	2.20	0.41
1:C:185:ARG:HD3	3:C:3521:HOH:O	2.19	0.41
1:A:47:LYS:HB2	1:A:47:LYS:HE3	1.85	0.41
1:B:61:PRO:HB2	1:B:64:VAL:HG23	2.03	0.41
1:C:153:LEU:C	1:C:153:LEU:HD22	2.40	0.41
1:C:105:ARG:HG2	3:C:3490:HOH:O	2.20	0.41
1:B:183:ARG:HB3	1:B:183:ARG:NH1	2.36	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	182/190 (96%)	180 (99%)	2 (1%)	0	100	100
1	B	172/190 (90%)	170 (99%)	2 (1%)	0	100	100
1	C	175/190 (92%)	173 (99%)	2 (1%)	0	100	100
All	All	529/570 (93%)	523 (99%)	6 (1%)	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	164/165 (99%)	157 (96%)	7 (4%)	40	16
1	B	158/165 (96%)	155 (98%)	3 (2%)	69	50

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	160/165 (97%)	157 (98%)	3 (2%)	69	50
All	All	482/495 (97%)	469 (97%)	13 (3%)	57	33

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	ARG
1	A	34	VAL
1	A	62	LYS
1	A	136	LYS
1	A	169	ASP
1	A	184	GLU
1	A	185	ARG
1	B	58	MSE
1	B	132	LEU
1	B	169	ASP
1	C	63	GLU
1	C	153	LEU
1	C	185	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	170	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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

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$
2	GDP	A	1401	-	30,30,30	1.66	7 (23%)	44,47,47	2.16	12 (27%)
2	GDP	B	2401	-	30,30,30	1.55	6 (20%)	44,47,47	2.75	11 (25%)
2	GDP	C	3401	-	30,30,30	1.33	5 (16%)	44,47,47	2.43	10 (22%)

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	GDP	A	1401	-	-	0/16/32/32	0/1/3/3
2	GDP	B	2401	-	-	0/16/32/32	0/1/3/3
2	GDP	C	3401	-	-	0/16/32/32	0/1/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1401	GDP	O4'-C1'	4.24	1.47	1.41
2	B	2401	GDP	C2-N2	3.86	1.38	1.32
2	A	1401	GDP	C2-N2	3.63	1.38	1.32
2	A	1401	GDP	C8-N9	3.44	1.41	1.36
2	B	2401	GDP	C8-N9	3.24	1.41	1.36
2	C	3401	GDP	C4-N3	-2.95	1.30	1.35
2	B	2401	GDP	C8-N7	-2.88	1.28	1.34
2	C	3401	GDP	C4-N9	-2.75	1.33	1.37
2	A	1401	GDP	C8-N7	-2.67	1.29	1.34
2	B	2401	GDP	C4-N9	-2.56	1.34	1.37
2	B	2401	GDP	C2-N3	-2.50	1.30	1.33
2	A	1401	GDP	O6-C6	-2.50	1.19	1.24
2	C	3401	GDP	C8-N7	-2.38	1.29	1.34
2	C	3401	GDP	C6-C5	2.20	1.44	1.41

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	3401	GDP	C2-N2	2.19	1.35	1.32
2	A	1401	GDP	O2'-C2'	2.16	1.48	1.43
2	A	1401	GDP	O4'-C4'	2.13	1.50	1.45
2	B	2401	GDP	C2-N1	2.03	1.40	1.36

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2401	GDP	C6-C5-N7	13.78	136.00	134.14
2	C	3401	GDP	C6-C5-N7	11.24	135.65	134.14
2	A	1401	GDP	C6-C5-N7	8.70	135.31	134.14
2	A	1401	GDP	O3'-C3'-C4'	4.59	124.61	111.08
2	B	2401	GDP	N3-C4-N9	4.35	133.29	126.91
2	B	2401	GDP	O3'-C3'-C4'	4.31	123.78	111.08
2	C	3401	GDP	O3'-C3'-C4'	4.28	123.69	111.08
2	A	1401	GDP	C2-N3-C4	4.12	120.88	115.09
2	A	1401	GDP	N3-C4-N9	3.93	132.67	126.91
2	C	3401	GDP	N3-C4-N9	3.83	132.52	126.91
2	B	2401	GDP	C2-N3-C4	3.71	120.30	115.09
2	C	3401	GDP	C2-N3-C4	3.64	120.21	115.09
2	C	3401	GDP	C3'-C2'-C1'	-3.57	95.32	100.91
2	A	1401	GDP	C5-C4-N3	-3.54	120.81	125.94
2	C	3401	GDP	C4-C5-N7	-3.48	106.55	109.52
2	B	2401	GDP	C5-C4-N3	-3.39	121.02	125.94
2	B	2401	GDP	C3'-C2'-C1'	-3.36	95.65	100.91
2	C	3401	GDP	C5-C4-N3	-3.19	121.33	125.94
2	B	2401	GDP	C4-C5-N7	-3.15	106.82	109.52
2	C	3401	GDP	O3B-PB-O2B	2.72	118.22	107.61
2	B	2401	GDP	O3B-PB-O2B	2.57	117.62	107.61
2	B	2401	GDP	O2'-C2'-C1'	-2.56	103.49	111.23
2	C	3401	GDP	O2'-C2'-C1'	-2.55	103.50	111.23
2	A	1401	GDP	O3B-PB-O2B	2.51	117.36	107.61
2	A	1401	GDP	O4'-C1'-C2'	-2.47	102.98	106.77
2	C	3401	GDP	O3B-PB-O1B	2.36	118.16	110.44
2	A	1401	GDP	O3B-PB-O1B	2.29	117.93	110.44
2	A	1401	GDP	O2'-C2'-C1'	-2.26	104.39	111.23
2	A	1401	GDP	C4-C5-N7	-2.19	107.65	109.52
2	B	2401	GDP	C8-N9-C4	2.16	108.55	106.90
2	A	1401	GDP	C3'-C2'-C1'	-2.05	97.69	100.91
2	A	1401	GDP	PA-O3A-PB	2.04	137.68	131.68
2	B	2401	GDP	C8-N7-C5	2.02	109.84	103.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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 ⓘ

### 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, 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	184/190 (96%)	-0.25	2 (1%) 77 83	7, 15, 31, 43	0
1	B	176/190 (92%)	-0.35	2 (1%) 77 83	6, 13, 29, 40	0
1	C	179/190 (94%)	-0.13	7 (3%) 37 42	8, 15, 39, 44	0
All	All	539/570 (94%)	-0.25	11 (2%) 62 67	6, 14, 37, 44	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	34	VAL	5.1
1	C	35	THR	4.5
1	C	33	GLY	3.9
1	A	30	LYS	3.6
1	A	29	GLY	2.9
1	C	36	ARG	2.6
1	C	26	VAL	2.5
1	B	26	VAL	2.4
1	B	36	ARG	2.4
1	C	62	LYS	2.3
1	C	56	PHE	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands

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, 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	GDP	A	1401	28/28	0.06	-0.75	10,12,15,17	0
2	GDP	C	3401	28/28	0.06	-1.08	8,10,12,13	0
2	GDP	B	2401	28/28	0.05	-1.17	7,8,10,11	0

## 6.5 Other polymers

There are no such residues in this entry.