



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:40 PM GMT

PDB ID : 4U03  
Title : Structure of the vibrio cholerae di-nucleotide cyclase (DncV) in complex with GTP and 5MTHFGLU2  
Authors : Zhu, D.Y.; Xiang, Y.  
Deposited on : 2014-07-11  
Resolution : 2.04 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](mailto: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

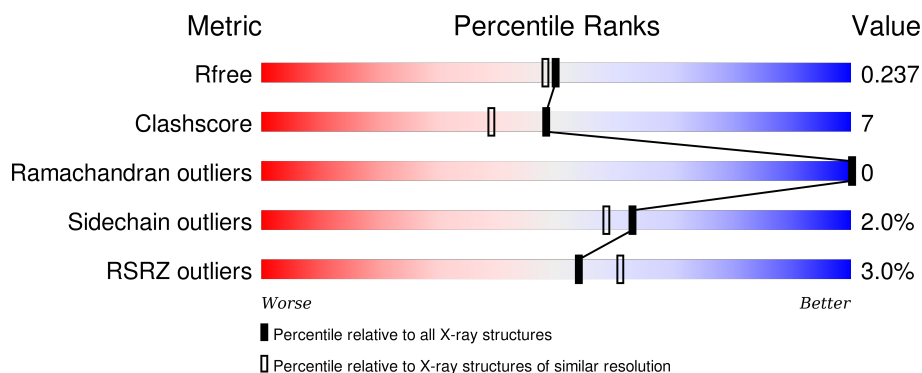
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.04 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	1192 (2.04-2.04)
Clashscore	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (2.04-2.04)

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. 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	427	<div> <div>2%</div> <div>79%</div> <div>11%</div> <div>9%</div> </div>
1	B	427	<div> <div>4%</div> <div>73%</div> <div>16%</div> <div>10%</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	GTP	A	502	-	-	-	X
2	GTP	B	502	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6803 atoms, of which 58 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.

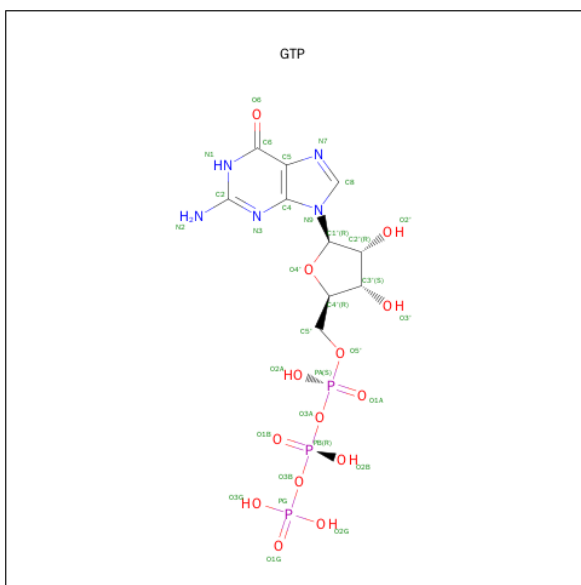
- Molecule 1 is a protein called Cyclic AMP-GMP synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	3	0
			3143	1986	542	599	16			
1	B	385	Total	C	N	O	S	0	2	0
			3101	1960	535	589	17			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	420	LEU	-	expression tag	UNP Q9KVG7
A	421	GLU	-	expression tag	UNP Q9KVG7
A	422	HIS	-	expression tag	UNP Q9KVG7
A	423	HIS	-	expression tag	UNP Q9KVG7
A	424	HIS	-	expression tag	UNP Q9KVG7
A	425	HIS	-	expression tag	UNP Q9KVG7
A	426	HIS	-	expression tag	UNP Q9KVG7
A	427	HIS	-	expression tag	UNP Q9KVG7
B	420	LEU	-	expression tag	UNP Q9KVG7
B	421	GLU	-	expression tag	UNP Q9KVG7
B	422	HIS	-	expression tag	UNP Q9KVG7
B	423	HIS	-	expression tag	UNP Q9KVG7
B	424	HIS	-	expression tag	UNP Q9KVG7
B	425	HIS	-	expression tag	UNP Q9KVG7
B	426	HIS	-	expression tag	UNP Q9KVG7
B	427	HIS	-	expression tag	UNP Q9KVG7

- Molecule 2 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula:  $C_{10}H_{16}N_5O_{14}P_3$ ).

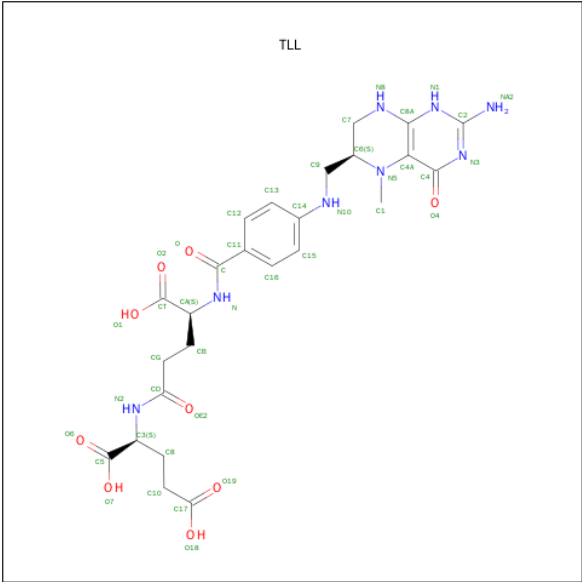


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	A	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	B	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	B	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Mg	0	0
			2	2		
3	A	2	Total	Mg	0	0
			2	2		

- Molecule 4 is N-[4-({[(6S)-2-amino-5-methyl-4-oxo-1,4,5,6,7,8-hexahydropteridin-6-yl]methyl}amino)benzoyl]-L-gamma-glutamyl-L-glutamic acid (three-letter code: TLL) (formula: C<sub>25</sub>H<sub>32</sub>N<sub>8</sub>O<sub>9</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	H	N	O	0	0
			71	25	29	8	9		
4	B	1	Total	C	H	N	O	0	0
			71	25	29	8	9		

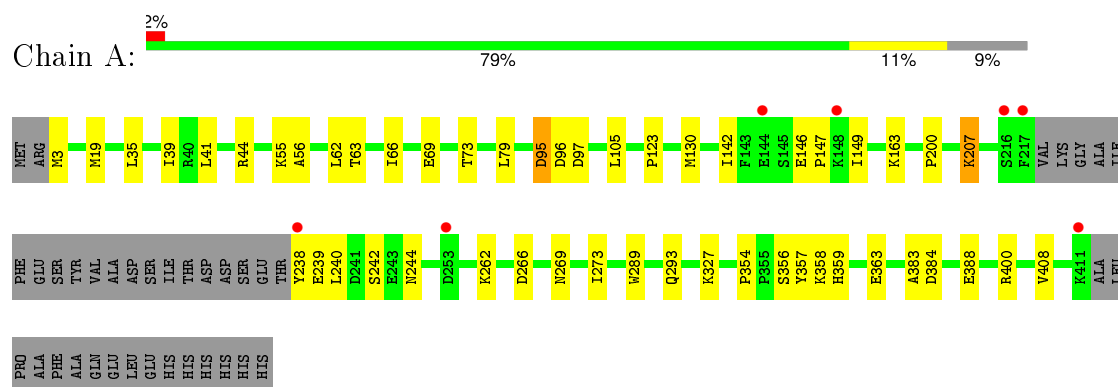
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	159	Total	O	0	0
			159	159		
5	B	126	Total	O	0	0
			126	126		

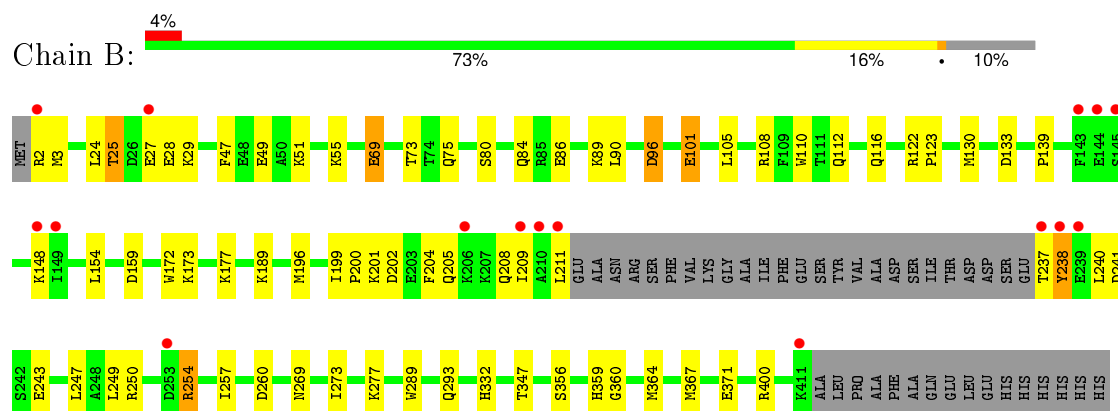
### 3 Residue-property plots [i](#)

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: Cyclic AMP-GMP synthase



#### • Molecule 1: Cyclic AMP-GMP synthase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	69.62Å 59.85Å 104.05Å 90.00° 95.81° 90.00°	Depositor
Resolution (Å)	43.63 – 2.04 43.63 – 2.04	Depositor EDS
% Data completeness (in resolution range)	99.5 (43.63-2.04) 95.7 (43.63-2.04)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.13 (at 2.05Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, $R_{free}$	0.180 , 0.233 0.186 , 0.237	Depositor DCC
$R_{free}$ test set	2755 reflections (5.56%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.5	Xtriage
Anisotropy	0.451	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 46.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 54236 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6803	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.81% of the height of the origin peak. No significant pseudotranslation is detected.*

<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, 0.375 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: GTP, MG, TLL

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	0.38	0/3214	0.52	1/4329 (0.0%)
1	B	0.37	0/3168	0.51	1/4268 (0.0%)
All	All	0.37	0/6382	0.51	2/8597 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	400	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	A	400	ARG	NE-CZ-NH2	-5.06	117.77	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	3143	0	3137	34	0
1	B	3101	0	3103	59	0
2	A	64	0	23	0	0
2	B	64	0	23	1	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	42	29	29	2	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	42	29	29	3	0
5	A	159	0	0	3	0
5	B	126	0	0	6	0
All	All	6745	58	6344	93	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 7.

All (93) 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:B:177:LYS:O	1:B:250:ARG:HD3	1.82	0.79
1:A:240:LEU:H	4:A:505:TLL:HN1	1.37	0.72
1:B:110:TRP:CH2	1:B:237:THR:HG21	2.25	0.71
1:B:189:LYS:HE3	5:B:723:HOH:O	1.94	0.68
1:B:211:LEU:HD22	1:B:238:TYR:HE2	1.63	0.64
1:B:172:TRP:O	1:B:173:LYS:HD2	1.99	0.63
1:B:172:TRP:C	1:B:173:LYS:HD2	2.19	0.63
1:A:239:GLU:HB2	1:A:262:LYS:HE2	1.81	0.62
1:A:357:TYR:CE1	1:A:358:LYS:HG3	2.36	0.61
1:A:383:ALA:HB1	1:A:388[A]:GLU:HG2	1.82	0.61
1:B:204:PHE:O	1:B:208:GLN:HG2	2.01	0.61
1:B:24:LEU:O	1:B:29:LYS:NZ	2.33	0.60
1:B:240:LEU:H	4:B:505:TLL:HN1	1.47	0.60
1:B:199:ILE:HB	1:B:200:PRO:HD2	1.84	0.60
1:A:44[B]:ARG:NH2	4:A:505:TLL:O1	2.27	0.59
1:B:25:THR:HB	1:B:27:GLU:HG2	1.83	0.59
1:A:289:TRP:O	1:A:293:GLN:HG2	2.02	0.58
1:A:262:LYS:HE3	1:A:266:ASP:OD2	2.04	0.57
1:A:95:ASP:OD1	1:A:95:ASP:N	2.30	0.56
1:B:247:LEU:HD23	1:B:257:ILE:HD11	1.88	0.56
1:B:86:GLU:O	1:B:90:LEU:HG	2.07	0.55
1:B:177:LYS:NZ	2:B:502:GTP:O3G	2.36	0.55
1:A:384:ASP:HB2	1:A:388[A]:GLU:OE1	2.05	0.55
1:B:28:GLU:OE2	1:B:122:ARG:NH2	2.39	0.55
1:A:123:PRO:HD3	1:A:130:MET:HG3	1.88	0.55
1:B:205:GLN:O	1:B:209:ILE:HG12	2.06	0.54
1:B:25:THR:OG1	1:B:28:GLU:OE2	2.26	0.53
1:B:2:ARG:HG3	1:B:3:MET:N	2.23	0.53
1:A:41:LEU:HD23	1:A:44[A]:ARG:HH21	1.73	0.53
1:B:123:PRO:HD3	1:B:130:MET:HG3	1.91	0.52

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:TYR:CD1	1:A:358:LYS:HG3	2.45	0.52
1:A:79:LEU:HD23	1:A:163:LYS:HD2	1.91	0.52
1:B:360:GLY:O	1:B:364:MET:HG2	2.10	0.51
1:A:269:ASN:O	1:A:273:ILE:HG12	2.11	0.51
1:B:204:PHE:HE1	4:B:505:TLL:C13	2.24	0.50
1:A:238:TYR:CG	1:A:239:GLU:N	2.79	0.50
1:B:211:LEU:HD13	1:B:238:TYR:CD2	2.47	0.50
1:B:356:SER:HA	1:B:359:HIS:CD2	2.47	0.50
1:A:19:MET:HG3	5:A:611:HOH:O	2.13	0.49
1:B:249:LEU:HD11	1:B:257:ILE:HG23	1.95	0.48
1:B:105:LEU:HD12	1:B:139:PRO:HD2	1.95	0.48
1:B:159:ASP:OD2	1:B:250:ARG:NH2	2.45	0.48
1:B:96:ASP:N	1:B:96:ASP:OD1	2.46	0.48
1:B:27:GLU:HG3	1:B:28:GLU:N	2.28	0.48
1:B:260:ASP:OD1	4:B:505:TLL:N3	2.46	0.48
1:B:154:LEU:HB3	1:B:196:MET:SD	2.54	0.48
1:B:241:ASP:OD1	1:B:243:GLU:HG2	2.14	0.47
1:A:207:LYS:HE3	1:A:239:GLU:O	2.14	0.47
1:B:249:LEU:HD11	1:B:257:ILE:CG2	2.43	0.47
1:B:277:LYS:HG2	5:B:615:HOH:O	2.13	0.47
1:B:332:HIS:HE1	5:B:617:HOH:O	1.98	0.47
1:A:142:ILE:HD11	1:A:200:PRO:HB3	1.97	0.47
1:A:142:ILE:HD11	1:A:200:PRO:HG3	1.96	0.46
1:B:69:GLU:O	1:B:73:THR:HG23	2.16	0.46
1:B:108:ARG:HG3	1:B:204:PHE:CZ	2.51	0.46
1:A:63:THR:OG1	1:A:66:ILE:HG13	2.16	0.46
1:B:101:GLU:OE2	1:B:148:LYS:HE3	2.16	0.46
1:B:201:LYS:O	1:B:205:GLN:HG3	2.16	0.46
1:B:25:THR:CB	1:B:27:GLU:HG2	2.47	0.45
1:B:80:SER:O	1:B:84:GLN:HG3	2.16	0.45
1:B:47:PHE:O	1:B:51:LYS:HG3	2.16	0.45
1:A:354:PRO:HG2	1:A:363:GLU:HG2	1.98	0.44
1:B:177:LYS:HE2	5:B:681:HOH:O	2.16	0.44
1:A:35:LEU:O	1:A:39:ILE:HD13	2.16	0.44
1:B:289:TRP:O	1:B:293:GLN:HG2	2.18	0.44
1:A:146:GLU:CD	1:A:147:PRO:HD2	2.38	0.44
1:B:27:GLU:HG3	1:B:28:GLU:H	1.82	0.44
1:A:56:ALA:HB1	1:A:62:LEU:HD11	1.99	0.44
1:B:116:GLN:NE2	5:B:726:HOH:O	2.51	0.43
1:B:367:MET:O	1:B:371:GLU:HG3	2.18	0.43
1:A:97:ASP:N	1:A:97:ASP:OD1	2.51	0.43

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:2:ARG:HB2	1:B:2:ARG:CZ	2.49	0.43
1:B:112:GLN:O	1:B:133:ASP:HB2	2.18	0.43
1:A:69:GLU:O	1:A:73:THR:HG23	2.18	0.43
1:B:332:HIS:HD2	5:B:619:HOH:O	2.02	0.42
1:B:249:LEU:CD1	1:B:257:ILE:HG23	2.48	0.42
1:A:356:SER:HA	1:A:359:HIS:CD2	2.54	0.42
1:B:254:ARG:NH1	1:B:257:ILE:HG22	2.35	0.42
1:B:208:GLN:O	1:B:211:LEU:HB3	2.21	0.41
1:A:55:LYS:HD2	5:A:713:HOH:O	2.21	0.41
1:B:159:ASP:CG	1:B:250:ARG:HH22	2.23	0.41
1:A:69:GLU:OE1	1:A:69:GLU:HA	2.21	0.41
1:B:55:LYS:HE3	1:B:55:LYS:HB2	1.65	0.41
1:A:356:SER:HA	1:A:359:HIS:CE1	2.55	0.41
1:B:89:LYS:HE2	1:B:89:LYS:HB3	1.87	0.41
1:A:44[A]:ARG:NE	5:A:685:HOH:O	2.35	0.41
1:B:27:GLU:CG	1:B:28:GLU:H	2.34	0.41
1:A:200:PRO:HG3	1:A:244:ASN:HB3	2.03	0.41
1:A:327:LYS:HB3	1:A:327:LYS:HE2	1.73	0.41
1:B:27:GLU:CG	1:B:28:GLU:N	2.84	0.40
1:B:269:ASN:O	1:B:273:ILE:HG23	2.21	0.40
1:B:49:GLU:OE2	1:B:75:GLN:HG3	2.22	0.40
1:A:105:LEU:HD22	1:A:149:ILE:HD13	2.03	0.40

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	388/427 (91%)	384 (99%)	4 (1%)	0	100	100
1	B	383/427 (90%)	377 (98%)	6 (2%)	0	100	100
All	All	771/854 (90%)	761 (99%)	10 (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	346/375 (92%)	340 (98%)	6 (2%)	68	65
1	B	342/375 (91%)	334 (98%)	8 (2%)	58	51
All	All	688/750 (92%)	674 (98%)	14 (2%)	63	58

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	MET
1	A	95	ASP
1	A	96	ASP
1	A	207	LYS
1	A	242	SER
1	A	408	VAL
1	B	25	THR
1	B	69	GLU
1	B	96	ASP
1	B	101	GLU
1	B	202	ASP
1	B	238	TYR
1	B	254	ARG
1	B	347	THR

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

Mol	Chain	Res	Type
1	B	208	GLN

### 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 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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$
2	GTP	A	501	3	25,34,34	1.12	2 (8%)	34,54,54	1.71	7 (20%)
2	GTP	A	502	3	25,34,34	1.27	3 (12%)	34,54,54	1.92	7 (20%)
4	TLL	A	505	-	30,44,44	4.00	13 (43%)	30,61,61	2.14	8 (26%)
2	GTP	B	501	3	25,34,34	1.13	3 (12%)	34,54,54	1.89	10 (29%)
2	GTP	B	502	3	25,34,34	1.21	3 (12%)	34,54,54	1.90	7 (20%)
4	TLL	B	505	-	30,44,44	4.26	16 (53%)	30,61,61	2.00	7 (23%)

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	GTP	A	501	3	-	0/18/38/38	0/3/3/3
2	GTP	A	502	3	-	0/18/38/38	0/3/3/3
4	TLL	A	505	-	-	0/25/48/48	0/3/3/3
2	GTP	B	501	3	-	0/18/38/38	0/3/3/3
2	GTP	B	502	3	-	0/18/38/38	0/3/3/3
4	TLL	B	505	-	-	0/25/48/48	0/3/3/3

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	505	TLL	C6-N5	-4.61	1.43	1.47
4	A	505	TLL	C6-N5	-4.27	1.43	1.47
4	A	505	TLL	CA-N	-2.52	1.43	1.46
4	A	505	TLL	C7-N8	-2.25	1.43	1.46
4	B	505	TLL	C13-C14	-2.23	1.35	1.39
4	A	505	TLL	C3-N2	-2.22	1.43	1.46
4	B	505	TLL	C12-C11	-2.09	1.35	1.39
4	B	505	TLL	C3-N2	-2.08	1.43	1.46
4	B	505	TLL	C7-N8	-2.03	1.43	1.46
2	B	502	GTP	O4'-C1'	2.05	1.43	1.41
4	B	505	TLL	CG-CD	2.07	1.55	1.51
2	B	501	GTP	O4'-C1'	2.19	1.44	1.41
4	B	505	TLL	C11-C	2.20	1.54	1.50
2	A	502	GTP	O4'-C1'	2.56	1.44	1.41
2	B	501	GTP	C5-C4	2.95	1.47	1.40
2	A	501	GTP	C5-C4	3.05	1.47	1.40
2	B	502	GTP	C5-C4	3.27	1.47	1.40
2	B	501	GTP	C6-C5	3.52	1.48	1.41
2	A	501	GTP	C6-C5	3.52	1.48	1.41
2	A	502	GTP	C5-C4	3.53	1.48	1.40
2	A	502	GTP	C6-C5	3.74	1.48	1.41
2	B	502	GTP	C6-C5	3.83	1.48	1.41
4	A	505	TLL	C-N	4.76	1.44	1.34
4	A	505	TLL	C8A-N8	4.86	1.41	1.35
4	A	505	TLL	CD-N2	4.99	1.44	1.34
4	B	505	TLL	C-N	5.09	1.45	1.34
4	B	505	TLL	C8A-N8	5.40	1.42	1.35
4	B	505	TLL	CD-N2	5.55	1.45	1.34
4	A	505	TLL	O4-C4	6.13	1.39	1.24
4	B	505	TLL	O4-C4	6.46	1.40	1.24
4	A	505	TLL	C12-C13	7.09	1.51	1.38
4	A	505	TLL	C2-NA2	7.47	1.43	1.32
4	A	505	TLL	C16-C11	7.72	1.52	1.39
4	A	505	TLL	C15-C14	7.92	1.52	1.39
4	B	505	TLL	C15-C14	8.02	1.52	1.39
4	B	505	TLL	C2-NA2	8.03	1.44	1.32
4	B	505	TLL	C12-C13	8.05	1.53	1.38
4	B	505	TLL	C16-C11	8.25	1.53	1.39
4	A	505	TLL	C4-N3	9.35	1.50	1.33
4	B	505	TLL	C4-N3	9.65	1.51	1.33

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	505	TLL	C7-C6-N5	-5.09	103.14	108.79
2	A	502	GTP	PA-O3A-PB	-4.64	119.69	132.73
2	B	502	GTP	C5-C6-N1	-4.56	117.35	123.59
2	A	502	GTP	C5-C6-N1	-4.52	117.41	123.59
4	B	505	TLL	C4A-C4-N3	-4.39	116.71	123.46
2	B	501	GTP	C5-C6-N1	-3.97	118.16	123.59
4	A	505	TLL	C4A-C4-N3	-3.91	117.44	123.46
4	B	505	TLL	N1-C2-N3	-3.78	116.73	121.79
2	A	501	GTP	C5-C6-N1	-3.77	118.43	123.59
2	B	501	GTP	C6-C5-C4	-3.76	116.41	120.90
2	B	502	GTP	C2'-C1'-N9	-3.73	108.60	114.29
2	B	501	GTP	N3-C2-N1	-3.48	122.14	127.44
2	A	502	GTP	PB-O3B-PG	-3.47	121.03	132.67
2	B	501	GTP	C2'-C1'-N9	-3.45	109.02	114.29
2	A	501	GTP	C6-C5-C4	-3.45	116.78	120.90
2	B	502	GTP	C4-C5-N7	-3.43	106.33	109.48
4	A	505	TLL	N1-C2-N3	-3.31	117.35	121.79
2	B	502	GTP	C6-C5-C4	-3.27	116.98	120.90
2	A	501	GTP	N3-C2-N1	-3.26	122.48	127.44
2	A	502	GTP	N3-C2-N1	-3.19	122.58	127.44
2	B	502	GTP	N3-C2-N1	-2.96	122.93	127.44
2	A	502	GTP	C4-C5-N7	-2.91	106.80	109.48
2	B	502	GTP	PA-O3A-PB	-2.68	125.20	132.73
2	A	501	GTP	C2'-C1'-N9	-2.67	110.21	114.29
2	A	501	GTP	C4-C5-N7	-2.67	107.02	109.48
2	A	502	GTP	C6-C5-C4	-2.66	117.72	120.90
4	A	505	TLL	OE2-CD-N2	-2.63	118.55	123.01
4	A	505	TLL	C10-C8-C3	-2.40	108.11	112.99
2	B	501	GTP	C4-C5-N7	-2.37	107.30	109.48
4	B	505	TLL	OE2-CD-N2	-2.31	119.08	123.01
2	B	501	GTP	C1'-N9-C4	-2.11	123.76	126.94
2	B	501	GTP	PA-O3A-PB	-2.05	126.99	132.73
2	A	501	GTP	O3G-PG-O2G	2.05	115.19	107.38
2	B	501	GTP	O2B-PB-O3B	2.12	114.69	105.09
2	B	501	GTP	O3G-PG-O2G	2.22	115.85	107.38
4	B	505	TLL	C4-N3-C2	2.33	119.49	116.06
4	B	505	TLL	CG-CD-N2	2.93	120.61	115.83
4	A	505	TLL	CG-CD-N2	3.03	120.78	115.83
4	A	505	TLL	NA2-C2-N1	3.66	121.80	117.80
2	A	501	GTP	C6-N1-C2	4.55	122.25	115.94
4	B	505	TLL	NA2-C2-N1	4.67	122.89	117.80
4	A	505	TLL	C4-C4A-C8A	4.85	118.29	114.43
2	B	502	GTP	C6-N1-C2	5.00	122.88	115.94

Continued on next page...



*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
2	B	501	GTP	C6-N1-C2	5.08	123.00	115.94
2	A	502	GTP	C6-N1-C2	5.13	123.06	115.94
4	B	505	TLL	C4-C4A-C8A	5.32	118.67	114.43

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	505	TLL	2	0
2	B	502	GTP	1	0
4	B	505	TLL	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, 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	389/427 (91%)	-0.20	7 (1%) 71 76	20, 35, 64, 91	0
1	B	385/427 (90%)	0.06	16 (4%) 40 46	21, 38, 77, 119	0
All	All	774/854 (90%)	-0.07	23 (2%) 54 61	20, 36, 72, 119	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	237	THR	6.5
1	B	209	ILE	5.9
1	B	211	LEU	5.9
1	B	238	TYR	5.8
1	B	210	ALA	4.7
1	B	145	SER	3.7
1	B	411	LYS	3.3
1	A	253	ASP	3.3
1	B	253	ASP	3.1
1	B	148	LYS	3.0
1	A	217	PHE	2.9
1	B	144	GLU	2.9
1	B	149	ILE	2.8
1	A	216	SER	2.8
1	B	206	LYS	2.6
1	B	27	GLU	2.4
1	A	144	GLU	2.4
1	A	411	LYS	2.3
1	B	2	ARG	2.3
1	B	143	PHE	2.2
1	A	148	LYS	2.2
1	A	238	TYR	2.1
1	B	239	GLU	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, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	GTP	A	502	32/32	0.78	0.17	2.38	48,71,147,149	0
2	GTP	B	502	32/32	0.88	0.13	2.21	34,54,126,134	0
4	TLL	B	505	42/42	0.83	0.22	0.55	48,72,104,110	0
2	GTP	A	501	32/32	0.98	0.11	0.40	24,37,45,51	0
4	TLL	A	505	42/42	0.94	0.12	-0.05	22,30,75,80	0
2	GTP	B	501	32/32	0.98	0.10	-0.68	28,38,48,50	0
3	MG	A	504	1/1	0.91	0.05	-	58,58,58,58	0
3	MG	B	504	1/1	0.92	0.14	-	52,52,52,52	0
3	MG	B	503	1/1	0.99	0.07	-	27,27,27,27	0
3	MG	A	503	1/1	0.98	0.06	-	26,26,26,26	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.