



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:45 AM UTC

PDB ID : 1CKM / pdb_00001ckm
Title : STRUCTURE OF TWO DIFFERENT CONFORMATIONS OF MRNA
CAPPING ENZYME IN COMPLEX WITH GTP
Authors : Hakansson, K.; Doherty, A.J.; Wigley, D.B.
Deposited on : 1997-04-20
Resolution : 2.50 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

<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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

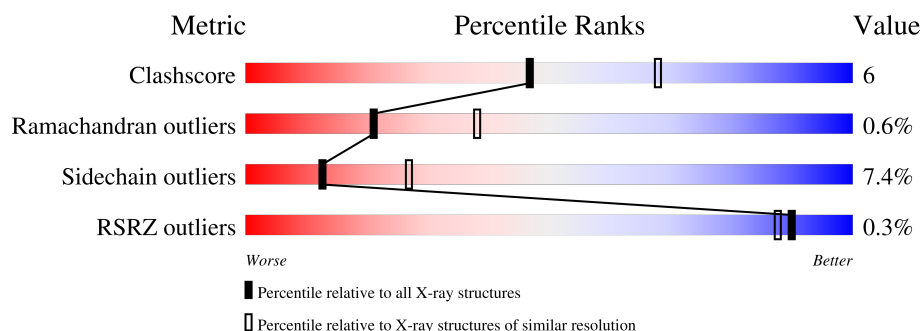
1 Overall quality at a glance

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))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	330	
1	B	330	

2 Entry composition [i](#)

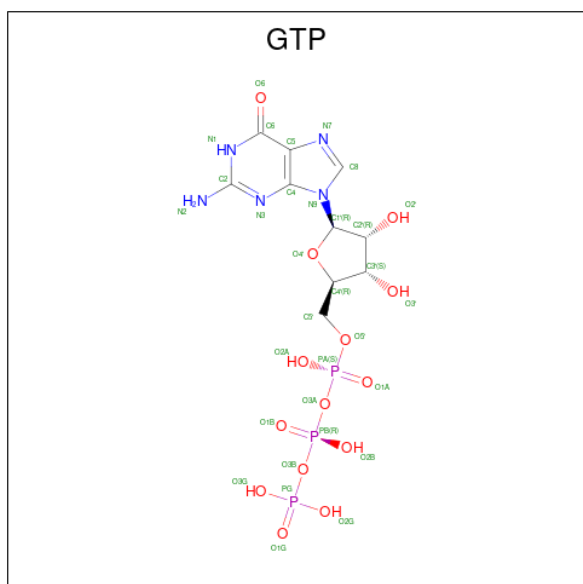
There are 3 unique types of molecules in this entry. The entry contains 5546 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 MRNA CAPPING ENZYME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	317	Total	C	N	O	S	0	0	0
			2561	1657	429	463	12			
1	B	317	Total	C	N	O	S	0	0	0
			2561	1657	429	463	12			

- Molecule 2 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: 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	B	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

- Molecule 3 is water.

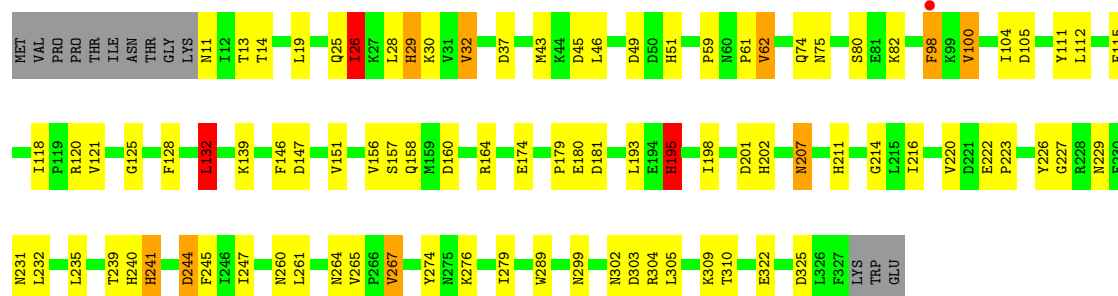
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	164	Total 164	O 164	0	0
3	B	196	Total 196	O 196	0	0

3 Residue-property plots [i](#)

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.

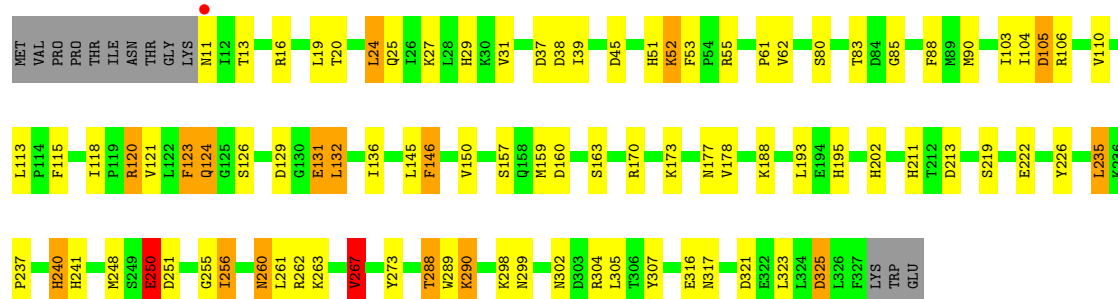
• Molecule 1: MRNA CAPPING ENZYME

Chain A: 



• Molecule 1: MRNA CAPPING ENZYME

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	93.33Å 214.93Å 105.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.50 10.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.5 (10.00-2.50) 93.1 (10.00-2.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.70 (at 2.50Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.218 , 0.299 0.199 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	34.2	Xtriage
Anisotropy	0.607	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 77.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5546	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% 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: GTP

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	0.94	7/2616 (0.3%)	1.66	43/3530 (1.2%)
1	B	0.96	8/2616 (0.3%)	1.71	51/3530 (1.4%)
All	All	0.95	15/5232 (0.3%)	1.69	94/7060 (1.3%)

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	A	0	3
1	B	0	3
All	All	0	6

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	211	HIS	CD2-NE2	-6.96	1.30	1.37
1	A	241	HIS	CD2-NE2	-6.87	1.30	1.37
1	B	211	HIS	CG-ND1	-6.37	1.31	1.38
1	B	195	HIS	CD2-NE2	-6.32	1.30	1.37
1	A	29	HIS	CD2-NE2	-6.28	1.30	1.37
1	B	51	HIS	CD2-NE2	-6.23	1.30	1.37
1	A	202	HIS	CD2-NE2	-6.21	1.31	1.37
1	A	240	HIS	CD2-NE2	-5.99	1.31	1.37
1	A	195	HIS	CD2-NE2	-5.96	1.31	1.37
1	A	51	HIS	CD2-NE2	-5.93	1.31	1.37
1	B	240	HIS	CD2-NE2	-5.84	1.31	1.37
1	B	241	HIS	CD2-NE2	-5.80	1.31	1.37
1	B	202	HIS	CD2-NE2	-5.77	1.31	1.37
1	B	29	HIS	CD2-NE2	-5.74	1.31	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	211	HIS	CD2-NE2	-5.34	1.31	1.37

All (94) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	207	ASN	OD1-CG-ND2	-9.49	113.11	122.60
1	A	105	ASP	CA-CB-CG	9.48	122.08	112.60
1	B	146	PHE	CA-CB-CG	9.39	123.19	113.80
1	A	98	PHE	CA-CB-CG	-9.00	104.80	113.80
1	B	195	HIS	CB-CG-CD2	-8.31	120.39	131.20
1	B	38	ASP	CA-CB-CG	8.31	120.91	112.60
1	A	157	SER	N-CA-C	8.21	121.15	111.71
1	B	105	ASP	CA-CB-CG	8.09	120.69	112.60
1	B	160	ASP	CA-CB-CG	7.96	120.56	112.60
1	A	267	VAL	N-CA-CB	-7.68	99.74	111.57
1	B	85	GLY	O-C-N	-7.65	117.33	123.14
1	B	288	THR	N-CA-CB	-7.60	99.65	111.05
1	A	160	ASP	CA-CB-CG	7.39	119.99	112.60
1	A	146	PHE	CA-CB-CG	7.27	121.07	113.80
1	A	303	ASP	CA-CB-CG	7.25	119.85	112.60
1	B	251	ASP	N-CA-C	7.25	121.70	113.01
1	A	231	ASN	OD1-CG-ND2	-7.24	115.36	122.60
1	A	325	ASP	CA-CB-CG	7.11	119.71	112.60
1	A	62	VAL	CB-CA-C	6.86	121.38	110.82
1	A	26	ILE	CB-CA-C	-6.84	99.97	110.50
1	B	302	ASN	OD1-CG-ND2	-6.67	115.92	122.60
1	B	316	GLU	CA-C-N	-6.55	114.02	122.79
1	B	316	GLU	C-N-CA	-6.55	114.02	122.79
1	A	49	ASP	CA-CB-CG	6.53	119.13	112.60
1	A	244	ASP	CA-CB-CG	6.49	119.09	112.60
1	B	123	PHE	CA-CB-CG	-6.48	107.32	113.80
1	B	211	HIS	CB-CG-CD2	-6.46	122.80	131.20
1	B	304	ARG	CA-CB-CG	6.44	126.97	114.10
1	B	302	ASN	CB-CG-ND2	6.41	126.02	116.40
1	A	195	HIS	CB-CG-CD2	-6.36	122.93	131.20
1	B	157	SER	N-CA-C	6.33	120.46	112.87
1	B	290	LYS	CB-CG-CD	-6.27	96.87	111.30
1	B	195	HIS	CB-CG-ND1	6.27	132.10	122.70
1	A	267	VAL	CB-CA-C	6.21	116.46	110.63
1	A	264	ASN	OD1-CG-ND2	-6.17	116.43	122.60
1	A	304	ARG	N-CA-C	6.17	117.67	111.07
1	B	11	ASN	CB-CG-ND2	6.15	125.62	116.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	55	ARG	CB-CA-C	-6.15	98.80	109.86
1	A	100	VAL	N-CA-CB	-6.13	102.57	111.52
1	A	211	HIS	CB-CG-CD2	-6.12	123.24	131.20
1	B	178	VAL	CA-C-N	6.08	127.43	119.84
1	B	178	VAL	C-N-CA	6.08	127.43	119.84
1	B	260	ASN	OD1-CG-ND2	-5.98	116.62	122.60
1	B	299	ASN	CA-CB-CG	5.97	118.57	112.60
1	B	195	HIS	CA-C-N	5.96	125.17	118.97
1	B	195	HIS	C-N-CA	5.96	125.17	118.97
1	A	45	ASP	CA-CB-CG	5.95	118.55	112.60
1	B	289	TRP	CE2-CD2-CG	-5.87	100.15	107.20
1	A	260	ASN	CA-CB-CG	5.84	118.44	112.60
1	A	261	LEU	N-CA-C	-5.81	105.29	112.90
1	B	129	ASP	CA-CB-CG	5.75	118.35	112.60
1	A	62	VAL	N-CA-CB	-5.71	100.09	111.91
1	B	250	GLU	CB-CG-CD	5.63	122.17	112.60
1	A	241	HIS	CB-CG-CD2	-5.56	123.97	131.20
1	B	124	GLN	OE1-CD-NE2	-5.55	117.05	122.60
1	B	51	HIS	CB-CG-CD2	-5.54	124.00	131.20
1	B	251	ASP	CA-CB-CG	5.52	118.12	112.60
1	B	53	PHE	CA-CB-CG	-5.52	108.28	113.80
1	B	11	ASN	CA-CB-CG	-5.49	107.11	112.60
1	B	120	ARG	CB-CG-CD	-5.48	98.69	111.30
1	B	256	ILE	CA-CB-CG2	5.42	119.71	110.50
1	A	51	HIS	CB-CG-CD2	-5.41	124.17	131.20
1	B	103	ILE	O-C-N	-5.36	117.25	122.93
1	A	265	VAL	CA-C-N	5.36	125.61	119.83
1	A	265	VAL	C-N-CA	5.36	125.61	119.83
1	B	317	ASN	OD1-CG-ND2	-5.31	117.29	122.60
1	B	289	TRP	CG-CD2-CE3	5.28	139.18	133.90
1	B	298	LYS	N-CA-C	-5.26	102.41	110.14
1	B	62	VAL	O-C-N	-5.24	117.53	123.03
1	A	11	ASN	CB-CG-ND2	5.22	124.23	116.40
1	A	37	ASP	CA-CB-CG	5.21	117.81	112.60
1	A	32	VAL	CB-CA-C	5.19	118.44	110.55
1	A	325	ASP	N-CA-C	-5.19	106.99	113.38
1	B	136	ILE	N-CA-CB	-5.19	103.98	110.47
1	B	131	GLU	CA-C-O	-5.19	114.72	120.38
1	A	62	VAL	CA-CB-CG1	5.17	119.19	110.40
1	A	289	TRP	CG-CD1-NE1	-5.17	103.48	110.20
1	B	29	HIS	CB-CG-CD2	-5.17	124.48	131.20
1	A	132	LEU	N-CA-C	-5.16	100.90	109.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	181	ASP	CA-CB-CG	5.15	117.75	112.60
1	B	213	ASP	CA-CB-CG	5.15	117.75	112.60
1	A	299	ASN	OD1-CG-ND2	-5.13	117.47	122.60
1	B	55	ARG	CA-CB-CG	5.12	124.35	114.10
1	A	226	TYR	N-CA-C	-5.11	100.84	109.07
1	B	260	ASN	CB-CG-ND2	5.10	124.05	116.40
1	B	177	ASN	CA-CB-CG	5.10	117.70	112.60
1	A	120	ARG	N-CA-C	5.09	118.59	112.38
1	B	267	VAL	N-CA-CB	-5.08	102.84	111.23
1	B	37	ASP	CA-CB-CG	5.08	117.68	112.60
1	B	202	HIS	CB-CG-CD2	-5.07	124.61	131.20
1	A	62	VAL	O-C-N	-5.06	117.72	123.03
1	A	202	HIS	CB-CG-CD2	-5.04	124.64	131.20
1	A	207	ASN	CB-CG-ND2	5.04	123.96	116.40
1	A	229	ASN	CA-CB-CG	5.04	117.64	112.60

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	111	TYR	Sidechain
1	A	274	TYR	Sidechain
1	A	98	PHE	Sidechain
1	B	226	TYR	Sidechain
1	B	273	TYR	Sidechain
1	B	307	TYR	Sidechain

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	2561	0	2607	27	0
1	B	2561	0	2607	30	0
2	A	32	0	12	3	0
2	B	32	0	12	3	0
3	A	164	0	0	1	0
3	B	196	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5546	0	5238	59	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 6.

All (59) 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:83:THR:HB	1:B:131:GLU:HG3	1.71	0.70
1:A:128:PHE:HE1	1:A:151:VAL:HG22	1.57	0.67
1:B:115:PHE:HB2	1:B:118:ILE:HD11	1.79	0.64
1:B:261:LEU:HB3	1:B:263:LYS:HG2	1.81	0.62
1:A:156:VAL:HG12	1:A:164:ARG:HG2	1.82	0.61
1:B:131:GLU:HG2	1:B:146:PHE:HZ	1.68	0.59
1:B:31:VAL:HG11	1:B:39:ILE:HD12	1.86	0.56
1:B:115:PHE:HB2	1:B:118:ILE:CD1	2.35	0.56
1:A:82:LYS:HE3	2:A:899:GTP:H5"	1.88	0.56
1:B:219:SER:OG	1:B:222:GLU:HB2	2.07	0.55
1:A:156:VAL:HG11	1:A:164:ARG:HA	1.89	0.54
1:B:321:ASP:O	1:B:325:ASP:HB2	2.08	0.53
1:A:26:ILE:HD11	1:A:132:LEU:HD11	1.89	0.53
1:A:195:HIS:HB3	1:A:198:ILE:HB	1.93	0.51
1:B:131:GLU:HG2	1:B:146:PHE:CZ	2.43	0.51
1:B:159:MET:CE	1:B:163:SER:HB3	2.41	0.51
1:A:115:PHE:HB2	1:A:118:ILE:HG12	1.91	0.50
1:B:240:HIS:NE2	1:B:290:LYS:HE2	2.26	0.50
1:B:121:VAL:HG23	1:B:124:GLN:OE1	2.12	0.49
1:A:158:GLN:HA	1:A:223:PRO:HB3	1.94	0.49
1:B:19:LEU:O	1:B:25:GLN:HA	2.14	0.48
1:B:170:ARG:O	1:B:173:LYS:HG2	2.14	0.47
1:A:214:GLY:HA3	1:A:235:LEU:O	2.14	0.47
1:B:88:PHE:CD2	1:B:132:LEU:HB2	2.49	0.47
1:B:159:MET:HE3	1:B:163:SER:HB3	1.97	0.47
1:A:59:PRO:HA	1:A:227:GLY:O	2.15	0.47
1:A:82:LYS:HG3	1:A:216:ILE:HD11	1.95	0.47
1:A:241:HIS:HB3	1:A:310:THR:HG23	1.97	0.47
1:B:145:LEU:O	1:B:188:LYS:HB2	2.15	0.46
1:A:164:ARG:NH2	1:A:222:GLU:O	2.48	0.46
1:A:305:LEU:HD23	3:A:1056:HOH:O	2.16	0.46
1:B:88:PHE:CE2	1:B:132:LEU:HB2	2.50	0.45
1:A:19:LEU:HB2	1:A:26:ILE:HG13	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:THR:HA	1:B:24:LEU:O	2.16	0.45
1:A:128:PHE:CE1	1:A:151:VAL:HG22	2.45	0.44
1:B:52:LYS:HD2	1:B:52:LYS:N	2.33	0.44
1:B:126:SER:HA	1:B:150:VAL:O	2.17	0.44
1:B:104:ILE:HG12	1:B:110:VAL:HG22	1.99	0.44
2:B:999:GTP:H4'	2:B:999:GTP:O2A	2.18	0.43
1:B:188:LYS:HE2	2:B:999:GTP:O6	2.18	0.43
1:A:74:GLN:HB2	1:A:75:ASN:ND2	2.34	0.43
1:A:147:ASP:OD1	1:A:164:ARG:HD2	2.19	0.42
1:B:120:ARG:HA	1:B:123:PHE:HD2	1.84	0.42
1:B:235:LEU:HD22	1:B:237:PRO:HD3	2.01	0.42
1:B:248:MET:SD	1:B:255:GLY:HA3	2.60	0.42
1:A:61:PRO:CG	2:A:899:GTP:H5'	2.50	0.42
1:A:244:ASP:H	1:A:302:ASN:ND2	2.18	0.42
1:A:245:PHE:O	1:A:279:ILE:HA	2.20	0.42
1:A:309:LYS:HD2	1:A:309:LYS:HA	1.87	0.42
1:A:26:ILE:HD12	1:A:28:LEU:HD11	2.01	0.41
1:B:90:MET:HE2	1:B:113:LEU:HD23	2.03	0.41
1:A:19:LEU:O	1:A:25:GLN:HA	2.21	0.41
1:A:43:MET:SD	1:A:104:ILE:HD13	2.61	0.41
2:A:899:GTP:N3	2:A:899:GTP:C2'	2.84	0.41
1:A:247:ILE:HG22	1:A:276:LYS:HA	2.02	0.40
1:B:61:PRO:HD2	2:B:999:GTP:H5''	2.02	0.40
1:B:80:SER:OG	1:B:188:LYS:HE3	2.21	0.40
1:B:105:ASP:HB2	3:B:1139:HOH:O	2.20	0.40
1:A:14:THR:HB	1:A:29:HIS:HB3	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	315/330 (96%)	301 (96%)	12 (4%)	2 (1%)	21	38
1	B	315/330 (96%)	303 (96%)	10 (3%)	2 (1%)	21	38
All	All	630/660 (96%)	604 (96%)	22 (4%)	4 (1%)	21	38

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	250	GLU
1	A	195	HIS
1	A	125	GLY
1	B	267	VAL

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	285/297 (96%)	262 (92%)	23 (8%)	11	23
1	B	285/297 (96%)	266 (93%)	19 (7%)	15	31
All	All	570/594 (96%)	528 (93%)	42 (7%)	13	27

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	THR
1	A	26	ILE
1	A	30	LYS
1	A	32	VAL
1	A	46	LEU
1	A	62	VAL
1	A	80	SER
1	A	100	VAL
1	A	112	LEU
1	A	121	VAL
1	A	132	LEU
1	A	139	LYS

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Mol	Chain	Res	Type
1	A	174	GLU
1	A	179	PRO
1	A	180	GLU
1	A	193	LEU
1	A	201	ASP
1	A	207	ASN
1	A	220	VAL
1	A	232	LEU
1	A	239	THR
1	A	267	VAL
1	A	322	GLU
1	B	13	THR
1	B	16	ARG
1	B	24	LEU
1	B	27	LYS
1	B	45	ASP
1	B	52	LYS
1	B	106	ARG
1	B	132	LEU
1	B	193	LEU
1	B	235	LEU
1	B	250	GLU
1	B	256	ILE
1	B	260	ASN
1	B	262	ARG
1	B	267	VAL
1	B	288	THR
1	B	305	LEU
1	B	323	LEU
1	B	325	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	29	HIS
1	A	75	ASN
1	A	117	ASN
1	A	177	ASN
1	A	211	HIS
1	A	264	ASN
1	A	317	ASN

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Mol	Chain	Res	Type
1	B	25	GLN
1	B	60	ASN
1	B	177	ASN
1	B	195	HIS
1	B	202	HIS
1	B	264	ASN
1	B	302	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	GTP	A	899	-	33,34,34	1.86	4 (12%)	50,54,54	0.76	0
2	GTP	B	999	-	33,34,34	1.70	8 (24%)	50,54,54	1.05	4 (8%)

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	899	-	-	3/22/38/38	0/3/3/3
2	GTP	B	999	-	-	1/22/38/38	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	899	GTP	PB-O3B	6.54	1.66	1.59
2	B	999	GTP	PA-O3A	5.83	1.65	1.59
2	A	899	GTP	PB-O3A	4.50	1.64	1.59
2	A	899	GTP	PA-O3A	4.46	1.64	1.59
2	B	999	GTP	PB-O3A	3.08	1.62	1.59
2	B	999	GTP	PB-O3B	-3.01	1.56	1.59
2	B	999	GTP	C2'-C3'	-2.78	1.45	1.53
2	B	999	GTP	C3'-C4'	-2.65	1.46	1.53
2	B	999	GTP	PG-O3G	-2.13	1.46	1.54
2	A	899	GTP	PG-O2G	-2.10	1.47	1.54
2	B	999	GTP	PG-O2G	-2.09	1.47	1.54
2	B	999	GTP	PA-O2A	-2.01	1.46	1.55

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	999	GTP	O4'-C4'-C5'	2.72	118.04	109.33
2	B	999	GTP	O4'-C1'-N9	2.15	113.22	108.36
2	B	999	GTP	O5'-C5'-C4'	2.14	116.27	108.99
2	B	999	GTP	O3G-PG-O2G	2.07	115.56	107.80

There are no chirality outliers.

All (4) torsion outliers are listed below:

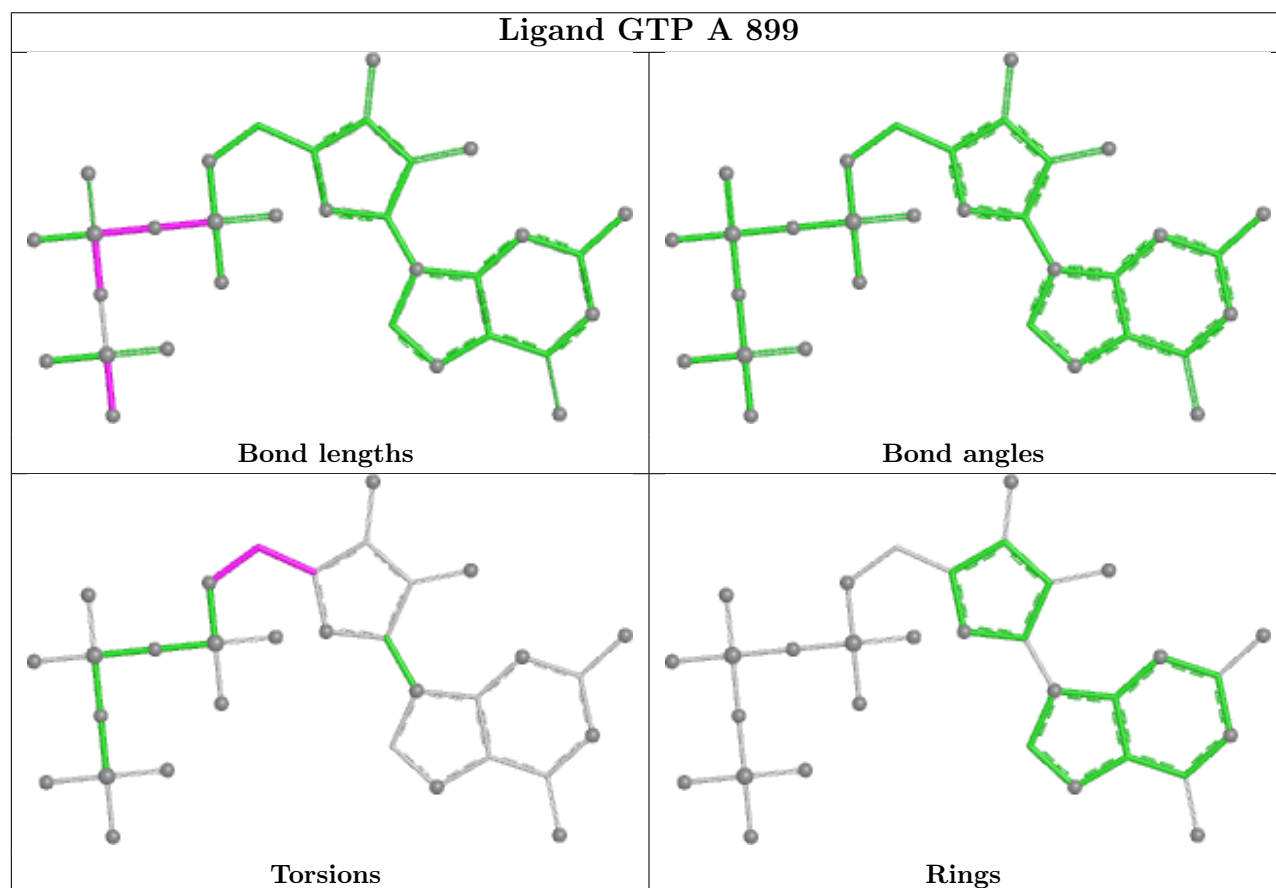
Mol	Chain	Res	Type	Atoms
2	B	999	GTP	C4'-C5'-O5'-PA
2	A	899	GTP	C3'-C4'-C5'-O5'
2	A	899	GTP	O4'-C4'-C5'-O5'
2	A	899	GTP	C4'-C5'-O5'-PA

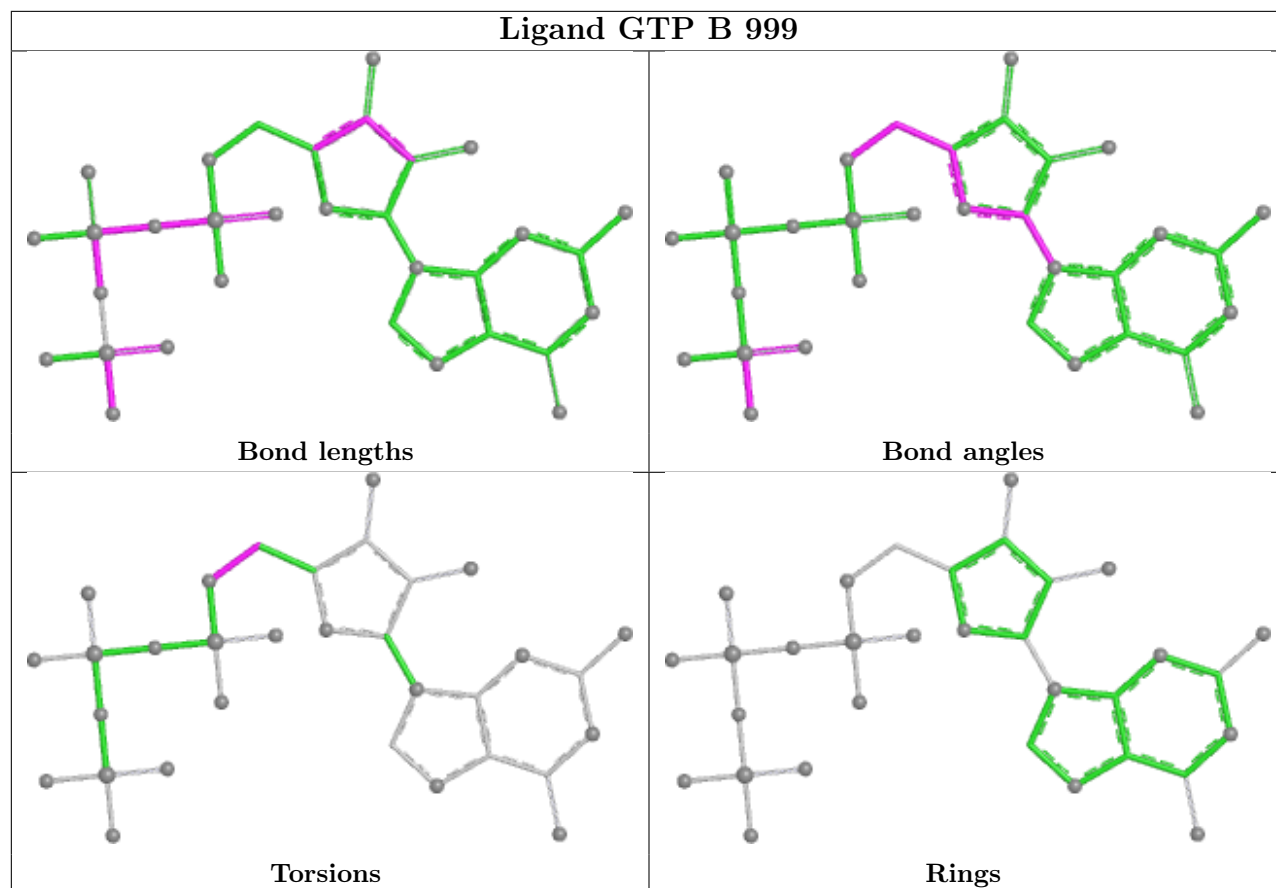
There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	899	GTP	3	0
2	B	999	GTP	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	317/330 (96%)	-0.61	1 (0%) 90 87	9, 24, 54, 71	0
1	B	317/330 (96%)	-0.60	1 (0%) 90 87	11, 24, 54, 93	0
All	All	634/660 (96%)	-0.60	2 (0%) 90 87	9, 24, 54, 93	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	11	ASN	2.4
1	A	98	PHE	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](#)

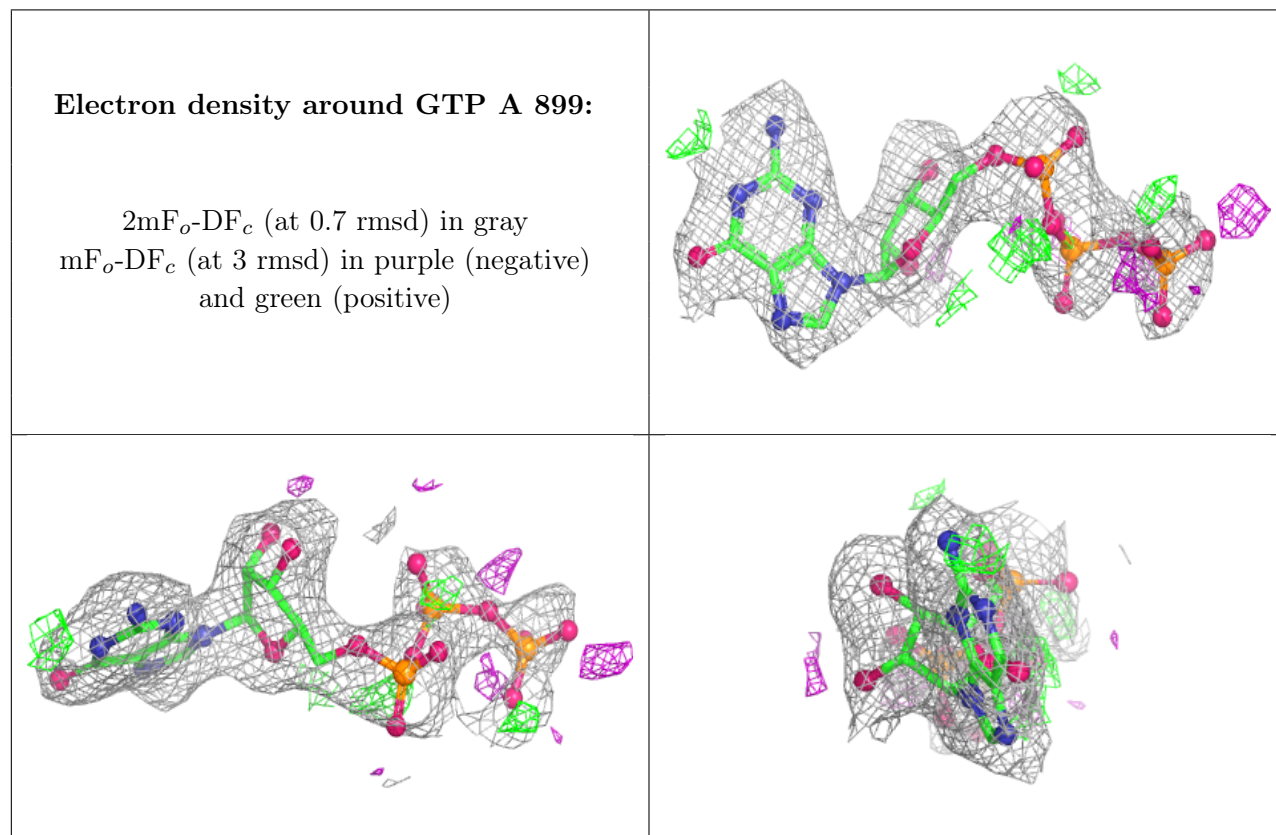
There are no oligosaccharides in this entry.

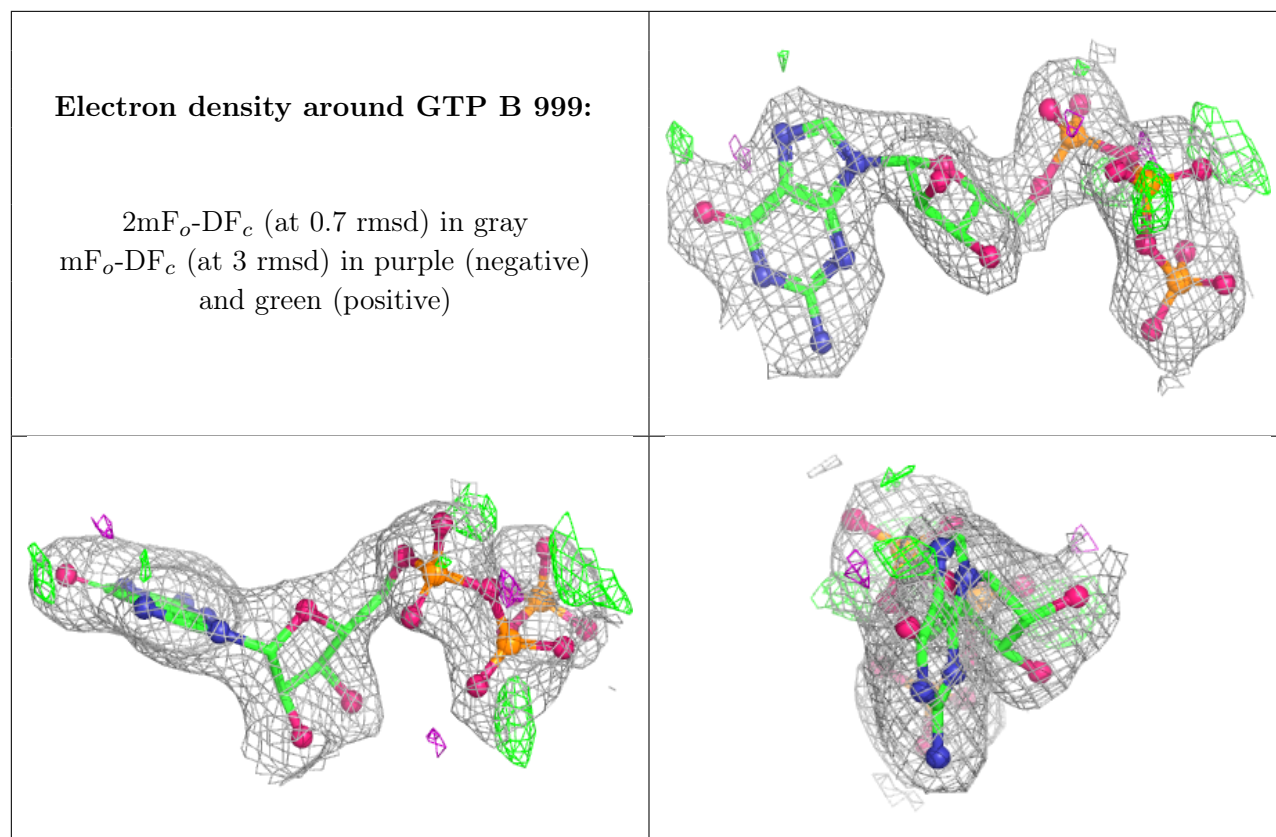
6.4 Ligands [i](#)

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, 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	B-factors(Å ²)	Q<0.9
2	GTP	A	899	32/32	0.91	0.10	16,35,124,125	0
2	GTP	B	999	32/32	0.97	0.06	10,19,42,50	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





6.5 Other polymers [i](#)

There are no such residues in this entry.