



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 02:54 AM GMT

PDB ID : 1CKN
Title : STRUCTURE OF GUANYLYLATED MRNA CAPPING ENZYME COM-
PLEXED 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 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/ValidationPDFNotes.html>

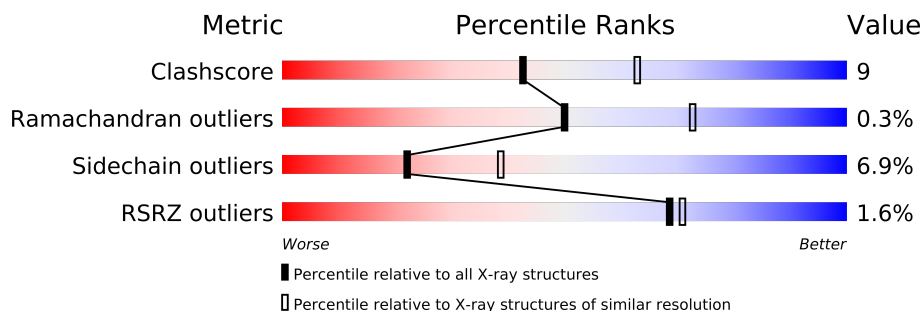
The following versions of software and data (see [references](#)) were used in the production of this report:

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 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	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (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. 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	330	
2	B	330	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	MN	B	1001	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5525 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 MRNA CAPPING ENZYME.

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

- Molecule 2 is a protein called MRNA CAPPING ENZYME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P S			
2	B	317	2584	1667	434	470	1 12	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	82	GPL	LYS	CONFLICT	UNP Q84424

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

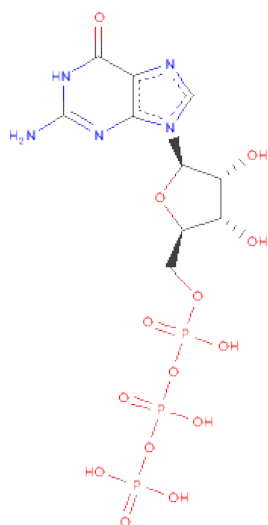
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Mn	0	0
			1	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		

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



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

- Molecule 6 is water.

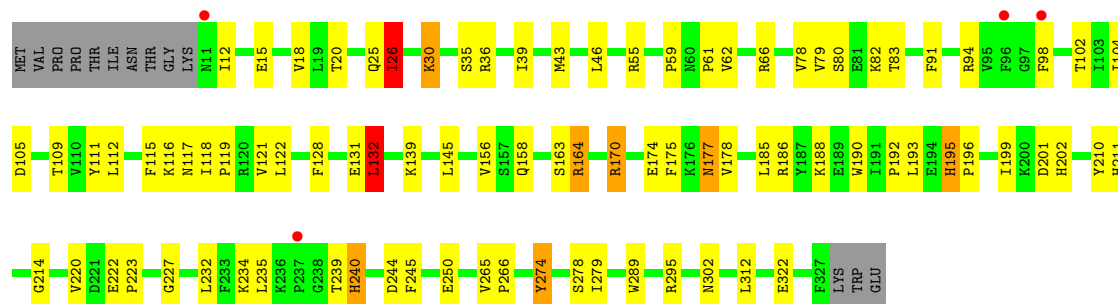
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	165	Total 165	O 165	0	0
6	B	177	Total 177	O 177	0	0

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: MRNA CAPPING ENZYME

Chain A: 



• Molecule 2: 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	94.91Å 212.95Å 105.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.50 14.99 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.6 (10.00-2.50) 95.4 (14.99-2.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.39 (at 2.51Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.224 , 0.299 0.206 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	35.1	Xtriage
Anisotropy	0.603	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 36093 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5525	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 4.76% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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: GPL, GTP, MN, SO4

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	0.78	0/2616	1.40	21/3530 (0.6%)
2	B	0.78	0/2606	1.45	33/3516 (0.9%)
All	All	0.78	0/5222	1.43	54/7046 (0.8%)

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	5
2	B	0	6
All	All	0	11

There are no bond length outliers.

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	186	ARG	NE-CZ-NH2	-10.63	114.98	120.30
2	B	289	TRP	CD1-CG-CD2	9.64	114.02	106.30
2	B	289	TRP	CE2-CD2-CG	-9.28	99.88	107.30
1	A	289	TRP	CD1-CG-CD2	9.22	113.67	106.30
1	A	164	ARG	NE-CZ-NH2	-8.76	115.92	120.30
2	B	66	ARG	NE-CZ-NH2	-8.46	116.07	120.30
2	B	164	ARG	NE-CZ-NH1	8.31	124.45	120.30
1	A	164	ARG	NE-CZ-NH1	8.03	124.32	120.30
1	A	289	TRP	CE2-CD2-CG	-7.91	100.97	107.30
2	B	164	ARG	NE-CZ-NH2	-7.83	116.39	120.30
2	B	106	ARG	NE-CZ-NH2	-7.82	116.39	120.30
2	B	106	ARG	NE-CZ-NH1	7.77	124.18	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	87	ARG	NE-CZ-NH1	7.49	124.05	120.30
1	A	186	ARG	NE-CZ-NH1	7.29	123.94	120.30
2	B	190	TRP	CD1-CG-CD2	7.23	112.08	106.30
1	A	170	ARG	NE-CZ-NH2	-7.22	116.69	120.30
2	B	190	TRP	CE2-CD2-CG	-7.05	101.66	107.30
2	B	87	ARG	NE-CZ-NH2	-6.98	116.81	120.30
2	B	304	ARG	NE-CZ-NH1	6.91	123.76	120.30
1	A	190	TRP	CE2-CD2-CG	-6.82	101.85	107.30
1	A	190	TRP	CD1-CG-CD2	6.57	111.56	106.30
2	B	120	ARG	NE-CZ-NH2	-6.32	117.14	120.30
2	B	170	ARG	NE-CZ-NH2	-6.32	117.14	120.30
2	B	55	ARG	CB-CA-C	-6.28	97.83	110.40
2	B	289	TRP	CG-CD2-CE3	6.17	139.45	133.90
2	B	267	VAL	N-CA-CB	-6.13	98.01	111.50
2	B	150	VAL	CG1-CB-CG2	-6.11	101.12	110.90
2	B	187	TYR	CB-CG-CD2	-6.08	117.35	121.00
2	B	55	ARG	NE-CZ-NH1	6.03	123.31	120.30
2	B	289	TRP	CB-CG-CD1	-6.00	119.20	127.00
1	A	94	ARG	NE-CZ-NH2	-5.91	117.34	120.30
2	B	228	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	A	240	HIS	CA-CB-CG	5.70	123.30	113.60
1	A	210	TYR	CB-CG-CD2	-5.70	117.58	121.00
2	B	55	ARG	CA-CB-CG	5.57	125.66	113.40
2	B	289	TRP	CG-CD1-NE1	-5.52	104.58	110.10
1	A	91	PHE	CB-CG-CD2	-5.51	116.94	120.80
2	B	218	MET	CG-SD-CE	-5.50	91.40	100.20
1	A	289	TRP	CG-CD1-NE1	-5.46	104.64	110.10
2	B	186	ARG	NE-CZ-NH2	-5.43	117.58	120.30
1	A	36	ARG	NE-CZ-NH1	5.42	123.01	120.30
2	B	295	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	A	274	TYR	CB-CG-CD2	-5.39	117.77	121.00
2	B	285	ALA	CB-CA-C	-5.33	102.10	110.10
1	A	132	LEU	CA-CB-CG	5.32	127.54	115.30
2	B	66	ARG	NE-CZ-NH1	5.29	122.94	120.30
2	B	251	ASP	CB-CG-OD1	5.26	123.04	118.30
1	A	26	ILE	CB-CA-C	-5.17	101.27	111.60
1	A	177	ASN	CB-CA-C	-5.16	100.08	110.40
2	B	304	ARG	CA-CB-CG	5.09	124.60	113.40
1	A	105	ASP	CA-CB-CG	5.09	124.59	113.40
2	B	295	ARG	NE-CZ-NH2	-5.07	117.77	120.30
1	A	66	ARG	NE-CZ-NH1	5.06	122.83	120.30
2	B	299	ASN	N-CA-CB	-5.01	101.58	110.60

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	111	TYR	Sidechain
1	A	170	ARG	Sidechain
1	A	274	TYR	Sidechain
1	A	295	ARG	Sidechain
1	A	98	PHE	Sidechain
2	B	106	ARG	Sidechain
2	B	111	TYR	Sidechain
2	B	187	TYR	Sidechain
2	B	273	TYR	Sidechain
2	B	274	TYR	Sidechain
2	B	77	TYR	Sidechain

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	2561	0	2607	41	0
2	B	2584	0	2617	54	0
3	B	1	0	0	0	0
4	B	5	0	0	0	0
5	A	32	0	12	5	0
6	A	165	0	0	7	0
6	B	177	0	0	4	0
All	All	5525	0	5236	93	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 9.

All (93) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:26:ILE:HD11	1:A:132:LEU:HD11	1.61	0.81
1:A:156:VAL:HG12	1:A:164:ARG:HG2	1.71	0.73
2:B:83:THR:HB	2:B:131:GLU:HG3	1.74	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:159:MET:CE	2:B:163:SER:HB3	2.26	0.66
1:A:158:GLN:HA	1:A:223:PRO:HB3	1.81	0.63
2:B:285:ALA:HB2	2:B:290:LYS:HD2	1.80	0.62
2:B:21:LEU:HD21	2:B:86:ILE:HG13	1.82	0.61
2:B:104:ILE:HG12	2:B:110:VAL:HG22	1.82	0.61
1:A:39:ILE:HD13	1:A:102:THR:HG21	1.82	0.61
2:B:202:HIS:HB2	6:B:1105:HOH:O	2.00	0.61
1:A:20:THR:HG23	1:A:25:GLN:HG2	1.84	0.60
1:A:61:PRO:HG2	5:A:899:GTP:H5'	1.82	0.60
2:B:131:GLU:HG2	2:B:146:PHE:HZ	1.68	0.58
2:B:248:MET:SD	2:B:255:GLY:HA3	2.43	0.58
2:B:77:TYR:HB2	2:B:193:LEU:HD12	1.85	0.58
2:B:100:VAL:HG12	2:B:102:THR:HG23	1.87	0.57
1:A:79:VAL:HG11	1:A:199:ILE:HD13	1.87	0.56
1:A:156:VAL:HG11	1:A:164:ARG:HA	1.87	0.56
2:B:145:LEU:O	2:B:188:LYS:HB2	2.06	0.55
1:A:214:GLY:HA3	1:A:235:LEU:O	2.07	0.55
2:B:321:ASP:O	2:B:325:ASP:HB2	2.06	0.55
1:A:18:VAL:HA	1:A:26:ILE:O	2.07	0.55
2:B:89:MET:HE1	2:B:129:ASP:HB2	1.88	0.55
1:A:12:ILE:HG23	1:A:35:SER:HB3	1.89	0.54
1:A:163:SER:HA	6:A:1026:HOH:O	2.08	0.54
2:B:159:MET:HE3	2:B:163:SER:HB3	1.88	0.54
2:B:159:MET:HE2	2:B:163:SER:HB3	1.88	0.54
2:B:82:GPL:HE2	2:B:234:LYS:HD2	1.89	0.54
2:B:95:VAL:HG12	2:B:96:PHE:CD2	2.43	0.54
2:B:61:PRO:HG3	2:B:232:LEU:HD13	1.88	0.54
1:A:82:LYS:HE3	5:A:899:GTP:H5''	1.90	0.53
2:B:121:VAL:HG22	2:B:152:SER:HB3	1.91	0.53
1:A:177:ASN:HD21	1:A:185:LEU:H	1.55	0.53
2:B:240:HIS:NE2	2:B:290:LYS:HE2	2.24	0.53
1:A:59:PRO:HA	1:A:227:GLY:O	2.09	0.52
2:B:115:PHE:HB2	2:B:118:ILE:CD1	2.40	0.52
2:B:121:VAL:HG23	2:B:124:GLN:OE1	2.09	0.52
1:A:119:PRO:HB2	1:A:122:LEU:HD23	1.92	0.51
2:B:64:ILE:HG21	2:B:235:LEU:HG	1.91	0.51
1:A:202:HIS:HB2	6:A:1012:HOH:O	2.11	0.50
2:B:244:ASP:H	2:B:302:ASN:ND2	2.09	0.50
1:A:15:GLU:O	1:A:30:LYS:N	2.44	0.50
1:A:188:LYS:NZ	5:A:899:GTP:O6	2.44	0.50
2:B:313:ASN:HA	6:B:1023:HOH:O	2.12	0.50
1:A:240:HIS:ND1	2:B:291:TYR:HB3	2.27	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:115:PHE:HB2	1:A:118:ILE:HG12	1.95	0.49
1:A:83:THR:HB	1:A:131:GLU:HG2	1.96	0.48
2:B:115:PHE:HE2	2:B:183:ALA:O	1.97	0.47
1:A:78:VAL:HG13	1:A:220:VAL:HG23	1.95	0.47
2:B:20:THR:HA	2:B:24:LEU:O	2.14	0.47
2:B:46:LEU:HD23	2:B:127:ILE:HG13	1.96	0.47
2:B:80:SER:OG	2:B:188:LYS:HE3	2.15	0.47
1:A:25:GLN:HG3	2:B:18:VAL:HG21	1.97	0.47
1:A:116:LYS:HD3	1:A:178:VAL:HG11	1.97	0.47
2:B:19:LEU:O	2:B:25:GLN:HA	2.15	0.46
1:A:61:PRO:CG	5:A:899:GTP:H5'	2.45	0.46
1:A:240:HIS:ND1	6:A:900:HOH:O	2.35	0.46
1:A:117:ASN:O	1:A:175:PHE:HD1	1.98	0.46
2:B:59:PRO:HB2	2:B:232:LEU:HD12	1.97	0.46
2:B:243:ILE:HA	2:B:302:ASN:HD22	1.79	0.46
1:A:30:LYS:HD2	1:A:109:THR:HG23	1.97	0.46
1:A:192:PRO:HB3	6:A:931:HOH:O	2.15	0.46
2:B:170:ARG:O	2:B:173:LYS:HG2	2.17	0.45
1:A:244:ASP:H	1:A:302:ASN:ND2	2.14	0.45
1:A:234:LYS:HD2	5:A:899:GTP:O2A	2.17	0.44
2:B:195:HIS:CE1	6:B:1107:HOH:O	2.70	0.44
2:B:92:PHE:HE2	2:B:118:ILE:HD13	1.83	0.44
2:B:14:THR:HG22	2:B:31:VAL:HA	1.99	0.44
1:A:211:HIS:HD2	6:A:1018:HOH:O	2.01	0.44
2:B:56:LEU:HD22	2:B:89:MET:HB2	2.01	0.43
2:B:115:PHE:CD1	2:B:118:ILE:HD11	2.53	0.43
2:B:85:GLY:HA3	2:B:132:LEU:O	2.19	0.43
2:B:88:PHE:CD2	2:B:132:LEU:HB2	2.54	0.43
1:A:128:PHE:HB3	1:A:145:LEU:HD22	2.01	0.42
1:A:195:HIS:CE1	6:A:1063:HOH:O	2.72	0.42
2:B:89:MET:CE	2:B:127:ILE:HG22	2.49	0.42
2:B:114:PRO:HD2	2:B:182:PRO:HG3	2.02	0.42
2:B:290:LYS:NZ	6:B:1174:HOH:O	2.53	0.42
2:B:92:PHE:CE2	2:B:118:ILE:HD13	2.53	0.42
2:B:279:ILE:N	2:B:279:ILE:HD12	2.35	0.41
2:B:49:ASP:OD2	2:B:55:ARG:HD2	2.19	0.41
2:B:285:ALA:CB	2:B:290:LYS:HD2	2.49	0.41
1:A:164:ARG:NH2	1:A:222:GLU:O	2.53	0.41
2:B:92:PHE:HD1	2:B:126:SER:HB2	1.86	0.41
2:B:131:GLU:HG2	2:B:146:PHE:CZ	2.53	0.41
1:A:196:PRO:HA	1:A:199:ILE:HD12	2.02	0.41
1:A:43:MET:SD	1:A:104:ILE:HD12	2.60	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:31:VAL:HG11	2:B:39:ILE:HD12	2.01	0.40
2:B:36:ARG:NH1	2:B:108:MET:O	2.54	0.40
1:A:278:SER:HB3	6:A:930:HOH:O	2.21	0.40
1:A:245:PHE:O	1:A:279:ILE:HA	2.22	0.40
2:B:203:LEU:O	2:B:207:ASN:ND2	2.50	0.40
1:A:265:VAL:HA	1:A:266:PRO:HD3	1.95	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	315/330 (96%)	298 (95%)	17 (5%)	0	100	100
2	B	314/330 (95%)	297 (95%)	15 (5%)	2 (1%)	33	55
All	All	629/660 (95%)	595 (95%)	32 (5%)	2 (0%)	50	73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	250	GLU
2	B	125	GLY

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%)	266 (93%)	19 (7%)	23	40
2	B	284/296 (96%)	264 (93%)	20 (7%)	21	38

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	569/593 (96%)	530 (93%)	39 (7%)	22	39

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	ILE
1	A	30	LYS
1	A	46	LEU
1	A	55	ARG
1	A	62	VAL
1	A	80	SER
1	A	112	LEU
1	A	121	VAL
1	A	132	LEU
1	A	139	LYS
1	A	174	GLU
1	A	193	LEU
1	A	195	HIS
1	A	201	ASP
1	A	232	LEU
1	A	239	THR
1	A	250	GLU
1	A	312	LEU
1	A	322	GLU
2	B	13	THR
2	B	16	ARG
2	B	27	LYS
2	B	45	ASP
2	B	52	LYS
2	B	55	ARG
2	B	132	LEU
2	B	189	GLU
2	B	193	LEU
2	B	232	LEU
2	B	235	LEU
2	B	250	GLU
2	B	256	ILE
2	B	262	ARG
2	B	267	VAL
2	B	288	THR
2	B	296	SER
2	B	305	LEU

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Mol	Chain	Res	Type
2	B	323	LEU
2	B	325	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	60	ASN
1	A	177	ASN
1	A	302	ASN
1	A	313	ASN
2	B	25	GLN
2	B	51	HIS
2	B	302	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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	GPL	B	82	3,2	34,34,35	3.67	10 (29%)	44,49,51	5.61	13 (29%)

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	GPL	B	82	3,2	-	1/18/37/39	0/1/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	82	GPL	O-C	18.66	1.24	1.11
2	B	82	GPL	P-O5'	5.56	1.62	1.57
2	B	82	GPL	P-O1P	4.00	1.51	1.46
2	B	82	GPL	CA-C	2.50	1.53	1.48
2	B	82	GPL	P-O2P	-2.46	1.48	1.55
2	B	82	GPL	C5-C4	-2.35	1.35	1.40
2	B	82	GPL	O4'-C1'	2.31	1.44	1.41
2	B	82	GPL	P-NZ	2.27	1.63	1.61
2	B	82	GPL	C8-N7	-2.25	1.30	1.34
2	B	82	GPL	C6-N1	2.24	1.41	1.37

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	82	GPL	C6-C5-N7	-32.28	129.79	134.14
2	B	82	GPL	P-NZ-CE	-9.86	111.40	124.67
2	B	82	GPL	O4'-C1'-N9	7.47	115.39	108.44
2	B	82	GPL	O1P-P-NZ	-6.75	100.82	113.71
2	B	82	GPL	C-CA-N	-6.64	107.19	113.83
2	B	82	GPL	O5'-P-NZ	4.41	120.09	106.24
2	B	82	GPL	C8-N9-C4	-3.28	104.39	106.90
2	B	82	GPL	C4'-O4'-C1'	-3.23	106.24	109.75
2	B	82	GPL	C6-N1-C2	2.87	124.53	119.51
2	B	82	GPL	O4'-C4'-C5'	2.81	119.39	109.36
2	B	82	GPL	O2P-P-O1P	2.65	116.02	109.89
2	B	82	GPL	O5'-C5'-C4'	2.60	118.49	108.94
2	B	82	GPL	O2P-P-NZ	-2.16	101.15	106.25

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	82	GPL	P-O5'-C5'-C4'

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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	$\# Z > 2$
5	GTP	A	899	-	34,34,34	1.52	7 (20%)	51,54,54	4.67	8 (15%)
4	SO4	B	1000	-	4,4,4	0.87	0	6,6,6	0.43	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
5	GTP	A	899	-	-	0/22/38/38	0/1/3/3
4	SO4	B	1000	-	-	0/0/0/0	0/0/0/0

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	899	GTP	O4'-C1'	3.29	1.46	1.41
5	A	899	GTP	PB-O3A	2.93	1.65	1.59
5	A	899	GTP	PB-O3B	2.77	1.64	1.59
5	A	899	GTP	PA-O2A	-2.72	1.42	1.55
5	A	899	GTP	C6-N1	2.22	1.40	1.37
5	A	899	GTP	PG-O3G	-2.14	1.46	1.54
5	A	899	GTP	PA-O3A	2.07	1.63	1.59

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	899	GTP	C6-C5-N7	-32.07	129.82	134.14
5	A	899	GTP	C6-N1-C2	3.92	126.37	119.51
5	A	899	GTP	C8-N9-C4	-3.02	104.59	106.90
5	A	899	GTP	C4'-O4'-C1'	-2.82	106.68	109.75
5	A	899	GTP	O4'-C1'-N9	-2.30	106.30	108.44
5	A	899	GTP	O5'-C5'-C4'	2.25	117.21	108.94
5	A	899	GTP	C4-C5-N7	2.09	111.31	109.52
5	A	899	GTP	O3G-PG-O2G	2.02	115.47	107.61

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, 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%)	-0.49	4 (1%) 74 76	8, 25, 55, 89	0
2	B	316/330 (95%)	-0.53	6 (1%) 64 66	7, 25, 56, 93	0
All	All	633/660 (95%)	-0.51	10 (1%) 68 71	7, 25, 55, 93	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	11	ASN	3.1
2	B	98	PHE	2.8
1	A	237	PRO	2.7
1	A	96	PHE	2.6
2	B	96	PHE	2.5
1	A	98	PHE	2.5
2	B	11	ASN	2.3
2	B	50	ASP	2.1
2	B	261	LEU	2.1
2	B	260	ASN	2.0

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

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	GPL	B	82	32/33	0.09	-0.30	4,16,25,31	0

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	MN	B	1001	1/1	0.13	6.21	28,28,28,28	0
5	GTP	A	899	32/32	0.14	0.70	15,36,130,135	0
4	SO4	B	1000	5/5	0.13	0.36	35,48,57,65	0

6.5 Other polymers

There are no such residues in this entry.