



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:48 PM GMT

PDB ID : 1CKO  
Title : STRUCTURE OF MRNA CAPPING ENZYME IN COMPLEX WITH THE  
CAP ANALOG GPPPG  
Authors : Hakansson, K.; Wigley, D.B.  
Deposited on : 1997-09-06  
Resolution : 3.10 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

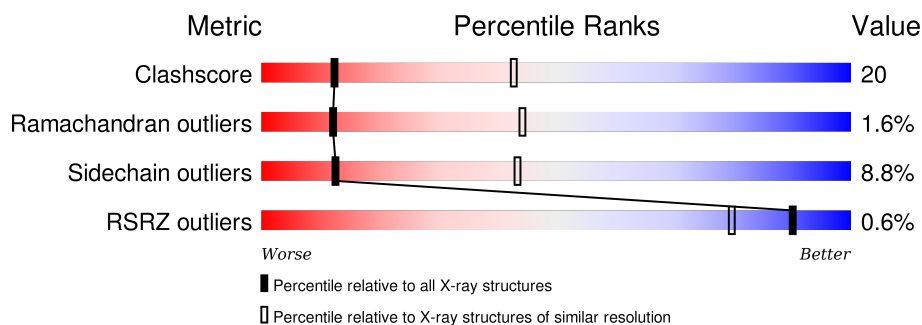
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.10 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	102246	1222 (3.14-3.06)
Ramachandran outliers	100387	1174 (3.14-3.06)
Sidechain outliers	100360	1174 (3.14-3.06)
RSRZ outliers	91569	1119 (3.14-3.06)

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	330	<div> <div></div> <div>49%</div> <div>37%</div> <div>8%</div> <div>• •</div> </div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2613 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.

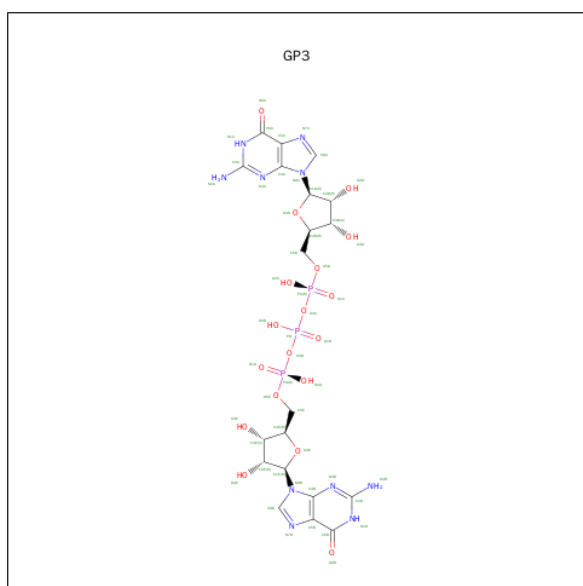
- Molecule 1 is a protein called MRNA CAPPING ENZYME.

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

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is DIGUANOSINE-5'-TRIPHOSPHATE (three-letter code: GP3) (formula:  $C_{20}H_{27}N_{10}O_{18}P_3$ ).

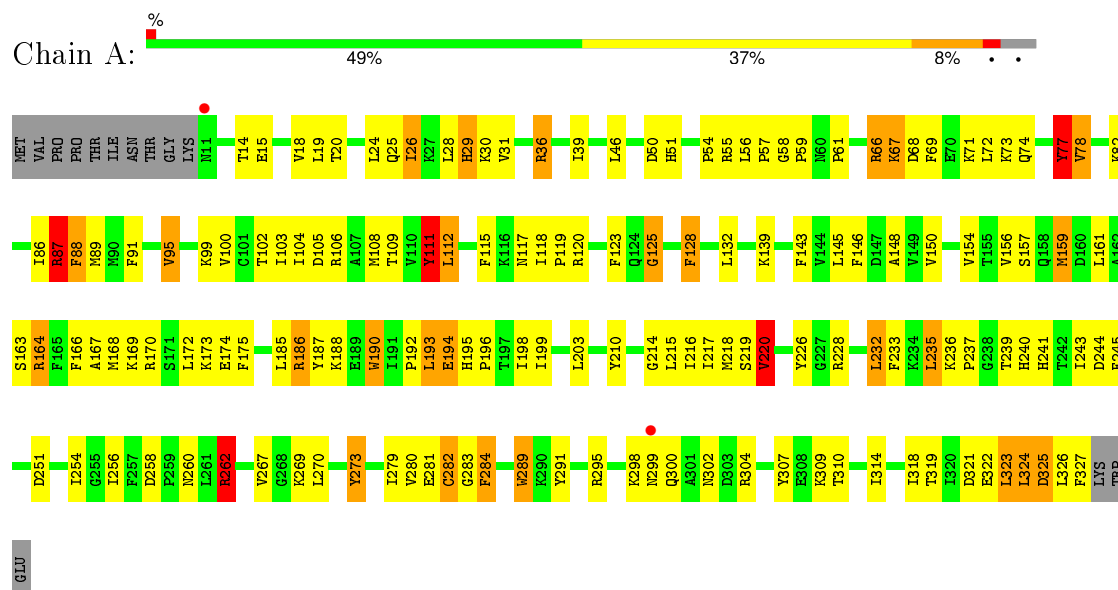


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			51	20	10	18	3		

### 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: MRNA CAPPING ENZYME



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	78.48Å 164.01Å 103.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 3.10 19.87 – 3.10	Depositor EDS
% Data completeness (in resolution range)	96.7 (10.00-3.10) 93.6 (19.87-3.10)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.40 (at 3.09Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.233 , 0.314 0.227 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	92.3	Xtriage
Anisotropy	0.299	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 119.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 12001 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	2613	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% 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: ZN, GP3

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	0.93	1/2616 (0.0%)	1.68	35/3530 (1.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	10

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	289	TRP	CG-CD2	-6.13	1.33	1.43

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	164	ARG	NE-CZ-NH2	9.40	125.00	120.30
1	A	120	ARG	NE-CZ-NH1	8.31	124.46	120.30
1	A	289	TRP	CD1-CG-CD2	8.26	112.91	106.30
1	A	120	ARG	NE-CZ-NH2	-8.14	116.23	120.30
1	A	190	TRP	CD1-CG-CD2	7.82	112.55	106.30
1	A	190	TRP	CE2-CD2-CG	-7.30	101.46	107.30
1	A	170	ARG	NE-CZ-NH2	-6.82	116.89	120.30
1	A	289	TRP	CE2-CD2-CG	-6.43	102.15	107.30
1	A	282	CYS	CA-CB-SG	-6.22	102.80	114.00
1	A	235	LEU	CA-C-N	-6.21	103.54	117.20
1	A	173	LYS	CA-CB-CG	6.17	126.97	113.40
1	A	66	ARG	CA-CB-CG	-6.16	99.84	113.40

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	146	PHE	CB-CA-C	-6.04	98.31	110.40
1	A	273	TYR	CB-CG-CD2	-5.93	117.44	121.00
1	A	324	LEU	CB-CG-CD1	-5.93	100.92	111.00
1	A	111	TYR	CB-CG-CD1	-5.88	117.47	121.00
1	A	106	ARG	NE-CZ-NH2	-5.87	117.36	120.30
1	A	128	PHE	CB-CG-CD2	-5.79	116.75	120.80
1	A	186	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	A	95	VAL	CG1-CB-CG2	5.46	119.63	110.90
1	A	220	VAL	CA-CB-CG2	-5.39	102.81	110.90
1	A	100	VAL	CA-CB-CG2	-5.39	102.81	110.90
1	A	66	ARG	NE-CZ-NH2	-5.39	117.61	120.30
1	A	170	ARG	NE-CZ-NH1	5.36	122.98	120.30
1	A	87	ARG	NE-CZ-NH2	-5.31	117.65	120.30
1	A	15	GLU	CA-CB-CG	5.30	125.06	113.40
1	A	146	PHE	CB-CG-CD2	-5.29	117.10	120.80
1	A	291	TYR	CB-CG-CD1	-5.27	117.83	121.00
1	A	106	ARG	NE-CZ-NH1	5.25	122.93	120.30
1	A	240	HIS	N-CA-CB	-5.24	101.17	110.60
1	A	210	TYR	CB-CA-C	-5.22	99.97	110.40
1	A	77	TYR	CB-CG-CD1	-5.17	117.90	121.00
1	A	300	GLN	CA-C-N	-5.16	105.86	117.20
1	A	88	PHE	CB-CG-CD1	-5.15	117.19	120.80
1	A	262	ARG	NE-CZ-NH1	5.05	122.83	120.30

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	111	TYR	Sidechain
1	A	187	TYR	Sidechain
1	A	226	TYR	Sidechain
1	A	273	TYR	Sidechain
1	A	284	PHE	Sidechain
1	A	304	ARG	Sidechain
1	A	307	TYR	Sidechain
1	A	77	TYR	Sidechain
1	A	87	ARG	Sidechain
1	A	91	PHE	Sidechain

## 5.2 Too-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 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	101	0
2	A	1	0	0	0	0
3	A	51	0	24	2	0
All	All	2613	0	2631	103	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 20.

All (103) 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:A:243:ILE:HG23	1:A:302:ASN:HD22	1.46	0.80
1:A:159:MET:SD	1:A:163:SER:HB3	2.27	0.74
1:A:279:ILE:HG21	1:A:298:LYS:HB2	1.71	0.72
1:A:77:TYR:HD1	1:A:219:SER:HA	1.58	0.69
1:A:77:TYR:CD1	1:A:219:SER:HA	2.30	0.66
1:A:281:GLU:HB2	1:A:295:ARG:HH11	1.61	0.65
1:A:217:ILE:HB	1:A:233:PHE:HB2	1.79	0.64
1:A:61:PRO:HG3	1:A:232:LEU:HD13	1.80	0.64
1:A:57:PRO:HG2	1:A:89:MET:HE3	1.81	0.62
1:A:244:ASP:H	1:A:302:ASN:ND2	1.98	0.62
1:A:57:PRO:O	1:A:89:MET:HE1	2.00	0.62
1:A:61:PRO:HA	1:A:232:LEU:HB3	1.81	0.61
1:A:145:LEU:HD11	1:A:185:LEU:HD13	1.81	0.61
1:A:128:PHE:HA	1:A:148:ALA:HA	1.83	0.61
1:A:56:LEU:HD22	1:A:89:MET:HB2	1.82	0.61
1:A:78:VAL:HG22	1:A:220:VAL:HG23	1.83	0.59
1:A:195:HIS:O	1:A:198:ILE:HB	2.02	0.59
1:A:192:PRO:HG2	1:A:195:HIS:CD2	2.38	0.59
1:A:156:VAL:HG12	1:A:164:ARG:HG2	1.84	0.58
1:A:72:LEU:HD23	1:A:193:LEU:HD11	1.84	0.58
1:A:73:LYS:HE3	1:A:326:LEU:HB3	1.84	0.58
1:A:82:LYS:HB2	1:A:216:ILE:HG12	1.85	0.58
1:A:161:LEU:HD12	1:A:164:ARG:NH1	2.19	0.58
1:A:323:LEU:HD13	1:A:324:LEU:HD23	1.86	0.57

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:HIS:HB3	1:A:314:ILE:HD11	1.87	0.56
1:A:19:LEU:HB2	1:A:26:ILE:HG13	1.87	0.54
1:A:194:GLU:HG2	1:A:195:HIS:CD2	2.42	0.54
1:A:194:GLU:HG2	1:A:195:HIS:HD2	1.74	0.53
1:A:20:THR:HA	1:A:24:LEU:O	2.09	0.53
1:A:14:THR:HG21	1:A:112:LEU:HD22	1.89	0.53
1:A:159:MET:HB3	1:A:163:SER:HB2	1.89	0.53
1:A:69:PHE:HA	1:A:72:LEU:HD13	1.91	0.53
1:A:196:PRO:HA	1:A:199:ILE:HD12	1.90	0.53
3:A:999:GP3:HO2B	3:A:999:GP3:HO3A	1.58	0.52
1:A:73:LYS:HA	1:A:327:PHE:HE2	1.74	0.51
1:A:86:ILE:HB	1:A:88:PHE:CE2	2.46	0.51
1:A:284:PHE:HD1	1:A:289:TRP:CE2	2.29	0.51
1:A:244:ASP:H	1:A:302:ASN:HD21	1.58	0.51
1:A:26:ILE:HD12	1:A:28:LEU:HD11	1.93	0.50
1:A:18:VAL:HA	1:A:26:ILE:O	2.11	0.50
1:A:283:GLY:O	1:A:289:TRP:HA	2.12	0.50
1:A:14:THR:HB	1:A:29:HIS:HB3	1.92	0.50
1:A:72:LEU:HA	1:A:77:TYR:HE2	1.76	0.50
1:A:39:ILE:HD13	1:A:102:THR:HG21	1.94	0.49
1:A:164:ARG:O	1:A:167:ALA:HB3	2.12	0.49
1:A:117:ASN:O	1:A:175:PHE:HD1	1.96	0.49
1:A:281:GLU:HB2	1:A:295:ARG:NH1	2.25	0.48
1:A:105:ASP:HB3	1:A:109:THR:HB	1.95	0.48
1:A:326:LEU:HB2	1:A:327:PHE:CD1	2.49	0.48
1:A:215:LEU:CD1	1:A:323:LEU:HD12	2.43	0.48
1:A:73:LYS:HA	1:A:327:PHE:CE2	2.48	0.48
1:A:66:ARG:HG3	1:A:318:ILE:HD11	1.96	0.48
1:A:172:LEU:O	1:A:175:PHE:HB3	2.14	0.48
1:A:87:ARG:HG3	1:A:88:PHE:N	2.28	0.48
1:A:269:LYS:HD3	1:A:270:LEU:O	2.14	0.48
1:A:190:TRP:CH2	1:A:218:MET:HE2	2.48	0.48
1:A:143:PHE:HB3	1:A:185:LEU:HD23	1.96	0.47
1:A:284:PHE:CD1	1:A:289:TRP:CE2	3.03	0.47
1:A:319:THR:OG1	1:A:322:GLU:HG3	2.14	0.47
1:A:166:PHE:CE1	1:A:169:LYS:HE3	2.50	0.46
1:A:67:LYS:HG3	1:A:68:ASP:N	2.29	0.46
1:A:256:ILE:HA	1:A:256:ILE:HD13	1.78	0.46
1:A:72:LEU:HA	1:A:77:TYR:CE2	2.50	0.46
1:A:258:ASP:O	1:A:262:ARG:N	2.46	0.46
1:A:284:PHE:HD1	1:A:289:TRP:CD2	2.34	0.45

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:LYS:HD2	1:A:123:PHE:CD2	2.52	0.45
1:A:125:GLY:O	1:A:150:VAL:HG12	2.16	0.45
1:A:243:ILE:HG23	1:A:302:ASN:ND2	2.23	0.45
1:A:26:ILE:HD11	1:A:132:LEU:HD21	1.98	0.45
1:A:14:THR:HA	1:A:30:LYS:O	2.17	0.45
1:A:71:LYS:HA	1:A:74:GLN:OE1	2.17	0.44
1:A:105:ASP:CB	1:A:109:THR:HB	2.47	0.44
1:A:128:PHE:HE2	1:A:172:LEU:HD11	1.83	0.44
1:A:148:ALA:HB2	1:A:168:MET:SD	2.56	0.44
1:A:103:ILE:C	1:A:104:ILE:HG13	2.38	0.44
1:A:251:ASP:O	1:A:269:LYS:HE2	2.17	0.44
1:A:326:LEU:HD12	1:A:327:PHE:HE1	1.83	0.44
1:A:51:HIS:CD2	1:A:55:ARG:HD3	2.53	0.43
1:A:109:THR:HG22	1:A:111:TYR:CE1	2.53	0.43
1:A:117:ASN:O	1:A:119:PRO:HD3	2.18	0.43
1:A:203:LEU:HD11	1:A:215:LEU:HD21	1.98	0.43
1:A:18:VAL:HG11	1:A:25:GLN:HG2	2.00	0.43
1:A:31:VAL:HG12	1:A:36:ARG:N	2.34	0.43
1:A:145:LEU:HB3	1:A:168:MET:CE	2.48	0.43
1:A:73:LYS:CE	1:A:326:LEU:HB3	2.48	0.43
1:A:86:ILE:HG22	1:A:87:ARG:O	2.18	0.43
1:A:216:ILE:HD13	1:A:216:ILE:HA	1.89	0.42
1:A:245:PHE:HB3	1:A:254:ILE:HB	2.01	0.42
1:A:194:GLU:C	1:A:196:PRO:HD2	2.40	0.42
1:A:161:LEU:O	1:A:164:ARG:HB2	2.19	0.42
1:A:310:THR:HG22	1:A:314:ILE:CD1	2.49	0.42
1:A:280:VAL:HG23	1:A:282:CYS:SG	2.59	0.42
1:A:54:PRO:HB2	1:A:108:MET:HG3	2.01	0.42
1:A:58:GLY:HA3	1:A:87:ARG:NH2	2.35	0.41
1:A:154:VAL:O	1:A:156:VAL:HG23	2.20	0.41
1:A:236:LYS:HA	1:A:237:PRO:HD3	1.78	0.41
1:A:118:ILE:HA	1:A:119:PRO:HD3	1.80	0.41
1:A:214:GLY:HA3	1:A:235:LEU:O	2.21	0.41
1:A:321:ASP:O	1:A:322:GLU:C	2.58	0.41
1:A:115:PHE:HB2	1:A:118:ILE:HG12	2.03	0.40
1:A:39:ILE:HG21	1:A:102:THR:HG21	2.03	0.40
3:A:999:GP3:O2E	3:A:999:GP3:O3D	2.33	0.40
1:A:59:PRO:HD2	1:A:87:ARG:NH2	2.36	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%)	284 (90%)	26 (8%)	5 (2%)	12	44

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	159	MET
1	A	67	LYS
1	A	194	GLU
1	A	325	ASP
1	A	125	GLY

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	285/297 (96%)	260 (91%)	25 (9%)	12	43

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	ILE
1	A	29	HIS
1	A	36	ARG
1	A	46	LEU
1	A	50	ASP
1	A	78	VAL

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	95	VAL
1	A	112	LEU
1	A	139	LYS
1	A	157	SER
1	A	174	GLU
1	A	186	ARG
1	A	188	LYS
1	A	193	LEU
1	A	220	VAL
1	A	228	ARG
1	A	232	LEU
1	A	239	THR
1	A	260	ASN
1	A	262	ARG
1	A	267	VAL
1	A	299	ASN
1	A	309	LYS
1	A	323	LEU
1	A	325	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	HIS
1	A	177	ASN
1	A	195	HIS
1	A	302	ASN
1	A	313	ASN

### 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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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$
3	GP3	A	999	-	42,56,56	1.56	8 (19%)	53,88,88	2.35	13 (24%)

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
3	GP3	A	999	-	-	0/24/64/64	0/6/6/6

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	999	GP3	C2D-C3D	-3.36	1.44	1.53
3	A	999	GP3	C2E-C3E	-2.75	1.45	1.53
3	A	999	GP3	C3D-C4D	-2.20	1.47	1.53
3	A	999	GP3	C8B-N7B	-2.20	1.30	1.34
3	A	999	GP3	C5E-C4E	2.54	1.59	1.51
3	A	999	GP3	C2B-N1B	2.67	1.40	1.35
3	A	999	GP3	C6A-N1A	3.29	1.39	1.33
3	A	999	GP3	C6B-N1B	4.49	1.41	1.33

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	999	GP3	C5B-C6B-N1B	-8.22	112.36	123.59
3	A	999	GP3	C5A-C6A-N1A	-7.18	113.77	123.59
3	A	999	GP3	N3B-C2B-N1B	-3.37	122.31	127.44
3	A	999	GP3	C4D-O4D-C1D	-3.15	106.26	109.72
3	A	999	GP3	N3A-C2A-N1A	-2.67	123.38	127.44
3	A	999	GP3	C5D-C4D-C3D	-2.18	106.55	115.21

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	A	999	GP3	O5E-C5E-C4E	2.22	117.31	109.12
3	A	999	GP3	C1D-N9A-C4A	2.27	130.36	126.94
3	A	999	GP3	C4A-C5A-N7A	2.69	111.95	109.48
3	A	999	GP3	O4E-C1E-N9B	2.93	114.22	108.10
3	A	999	GP3	C2D-C1D-N9A	3.62	119.83	114.29
3	A	999	GP3	C6A-N1A-C2A	5.55	123.65	115.94
3	A	999	GP3	C6B-N1B-C2B	5.98	124.24	115.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	999	GP3	2	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	317/330 (96%)	-0.57	2 (0%) 90 80	22, 66, 121, 144	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	11	ASN	2.6
1	A	299	ASN	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
3	GP3	A	999	51/51	0.96	0.15	-0.76	27,52,85,101	0
2	ZN	A	888	1/1	0.95	0.15	-	103,103,103,103	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.