



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 20, 2020 – 04:09 PM BST

PDB ID : 4HDG
Title : Crystal Structure of viral RdRp in complex with GTP
Authors : Surana, P.; Nair, D.T.
Deposited on : 2012-10-02
Resolution : 2.38 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

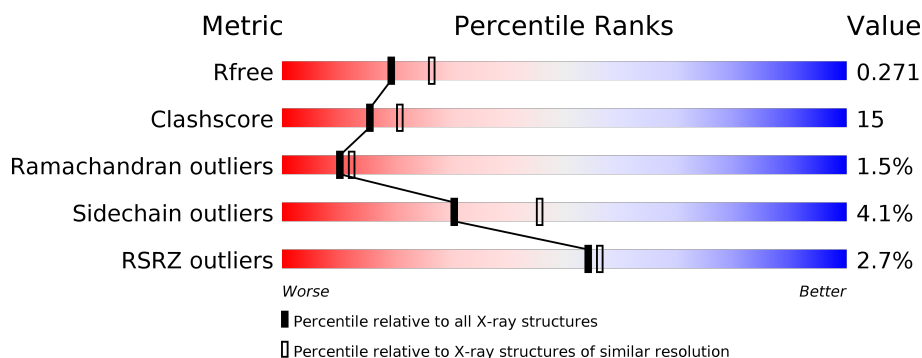
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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}	130704	5509 (2.40-2.36)
Clashscore	141614	6082 (2.40-2.36)
Ramachandran outliers	138981	5973 (2.40-2.36)
Sidechain outliers	138945	5975 (2.40-2.36)
RSRZ outliers	127900	5397 (2.40-2.36)

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 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	639	<div> <div>3%</div> <div> <div></div> <div>72%</div> <div>21%</div> <div>..</div> </div> </div>
1	B	639	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>17%</div> <div>...</div> </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	1001	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10361 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 Polypeptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	612	Total	C	N	O	S	0	0	0
			4924	3105	889	900	30			
1	B	612	Total	C	N	O	S	0	0	0
			4914	3100	884	899	31			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	267	GLY	-	EXPRESSION TAG	UNP G3LHD9
A	268	PRO	-	EXPRESSION TAG	UNP G3LHD9
A	269	LEU	-	EXPRESSION TAG	UNP G3LHD9
A	270	GLY	-	EXPRESSION TAG	UNP G3LHD9
A	271	SER	-	EXPRESSION TAG	UNP G3LHD9
A	373	ARG	LYS	CONFLICT	UNP G3LHD9
A	429	ASN	ASP	CONFLICT	UNP G3LHD9
A	836	ALA	THR	CONFLICT	UNP G3LHD9
B	267	GLY	-	EXPRESSION TAG	UNP G3LHD9
B	268	PRO	-	EXPRESSION TAG	UNP G3LHD9
B	269	LEU	-	EXPRESSION TAG	UNP G3LHD9
B	270	GLY	-	EXPRESSION TAG	UNP G3LHD9
B	271	SER	-	EXPRESSION TAG	UNP G3LHD9
B	373	ARG	LYS	CONFLICT	UNP G3LHD9
B	429	ASN	ASP	CONFLICT	UNP G3LHD9
B	836	ALA	THR	CONFLICT	UNP G3LHD9

- 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	B	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	211	Total	O	0	0
			211	211		
4	B	244	Total	O	0	0
			244	244		

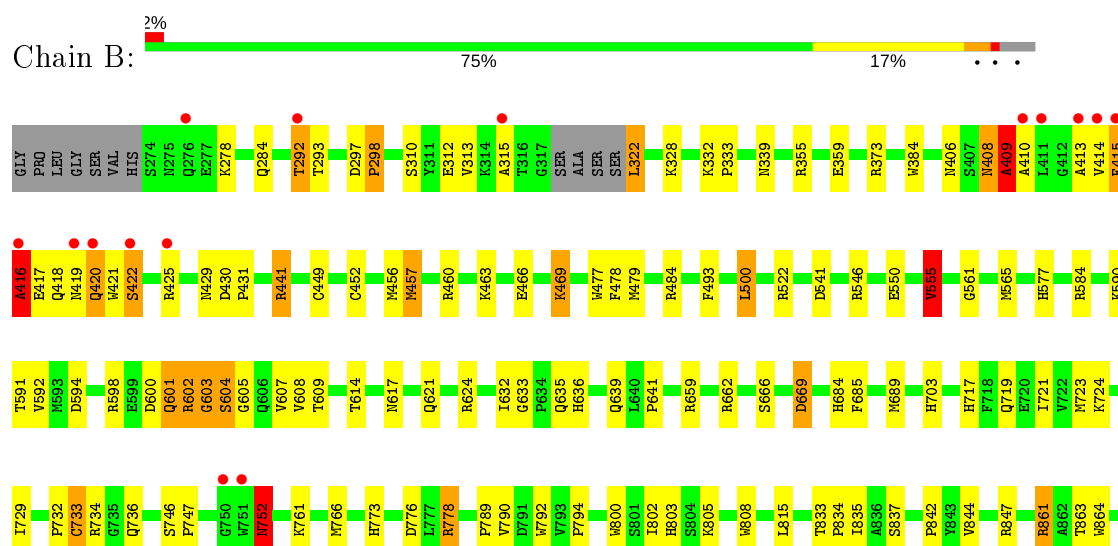
3 Residue-property plots [i](#)

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: Polyprotein



• Molecule 1: Polyprotein



I867	I868	I872	V885	T889	SER	LEU	ARG	ARG	TYR	GLU	ASP	VAL	LEU	ILE	GLN	GLU	ASP	ARG	VAL	ILE
------	------	------	------	------	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	144.57Å 86.55Å 112.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.38 49.77 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.7 (50.00-2.38) 99.8 (49.77-2.38)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.37Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.220 , 0.276 0.216 , 0.271	Depositor DCC
R_{free} test set	4034 reflections (7.04%)	wwPDB-VP
Wilson B-factor (Å ²)	32.0	Xtriage
Anisotropy	0.780	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10361	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 50.41 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.4982e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.85	3/5041 (0.1%)	0.82	6/6823 (0.1%)
1	B	0.83	2/5031 (0.0%)	0.81	6/6811 (0.1%)
All	All	0.84	5/10072 (0.0%)	0.82	12/13634 (0.1%)

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	2
1	B	0	2
All	All	0	4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	422	SER	CB-OG	10.26	1.55	1.42
1	B	733	CYS	CB-SG	-8.15	1.68	1.82
1	B	298	PRO	N-CD	7.44	1.58	1.47
1	A	416	ALA	CA-CB	7.16	1.67	1.52
1	A	834	PRO	N-CD	7.13	1.57	1.47

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	392	ARG	NE-CZ-NH2	-9.02	115.79	120.30
1	A	500	LEU	CA-CB-CG	-7.24	98.65	115.30
1	A	502	ARG	NE-CZ-NH2	-7.19	116.70	120.30
1	A	861	ARG	NE-CZ-NH1	6.25	123.42	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	409	ALA	N-CA-C	6.25	127.86	111.00
1	B	500	LEU	CA-CB-CG	-5.71	102.16	115.30
1	B	662	ARG	NE-CZ-NH2	-5.66	117.47	120.30
1	A	297	ASP	C-N-CD	5.65	140.26	128.40
1	B	624	ARG	NE-CZ-NH2	-5.25	117.67	120.30
1	B	555	VAL	CB-CA-C	-5.23	101.46	111.40
1	A	555	VAL	CB-CA-C	-5.17	101.58	111.40
1	B	778	ARG	NE-CZ-NH1	5.17	122.88	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	300	HIS	Mainchain
1	A	416	ALA	Peptide
1	B	408	ASN	Peptide
1	B	416	ALA	Peptide

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	4924	0	4826	156	0
1	B	4914	0	4808	131	0
2	A	32	0	12	9	0
2	B	32	0	12	2	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	211	0	0	7	0
4	B	244	0	0	15	0
All	All	10361	0	9658	289	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 15.

All (289) 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:527:LYS:NZ	1:A:661:THR:OG1	1.79	1.13
1:A:742:ARG:HH11	1:A:742:ARG:HG2	0.96	1.12
1:A:487:GLU:CD	1:A:607:VAL:HG12	1.70	1.11
1:A:416:ALA:HB3	1:A:478:PHE:CE2	1.86	1.10
1:A:414:VAL:HG11	1:A:478:PHE:CE1	1.86	1.09
1:B:284:GLN:HG3	4:B:1297:HOH:O	1.51	1.07
1:A:298:PRO:HB2	1:A:584:ARG:HH12	1.20	1.07
1:A:457:MET:HG3	1:A:478:PHE:CD1	1.89	1.06
1:A:602:ARG:HG2	1:A:603:GLY:H	1.21	1.04
1:A:330:MET:HE1	1:A:783:ALA:HB1	1.40	1.03
1:A:416:ALA:HB3	1:A:478:PHE:HE2	1.19	0.98
1:B:602:ARG:HG3	1:B:604:SER:H	1.27	0.97
1:A:742:ARG:NH1	1:A:742:ARG:HG2	1.74	0.97
1:A:633:GLY:H	1:A:636:HIS:HD2	1.13	0.94
1:A:603:GLY:HA3	4:A:1280:HOH:O	1.68	0.93
1:B:292:THR:HG22	1:B:293:THR:HG23	1.50	0.91
1:B:633:GLY:H	1:B:636:HIS:CD2	1.90	0.89
1:A:633:GLY:H	1:A:636:HIS:CD2	1.92	0.88
1:B:633:GLY:H	1:B:636:HIS:HD2	1.18	0.88
1:A:634:PRO:HD2	1:A:635:GLN:HE21	1.41	0.85
1:A:416:ALA:HA	1:A:418:GLN:HE21	1.41	0.84
1:A:414:VAL:HG11	1:A:478:PHE:CZ	2.12	0.83
1:B:766:MET:HG3	1:B:808:TRP:CE2	2.13	0.83
1:A:600:ASP:CG	1:A:601:GLN:H	1.81	0.83
1:B:600:ASP:CG	1:B:601:GLN:H	1.79	0.83
1:B:406:ASN:OD1	1:B:425:ARG:NH1	2.12	0.82
1:B:359:GLU:OE1	4:B:1217:HOH:O	1.98	0.81
1:B:790:VAL:HG21	1:B:885:VAL:HG23	1.64	0.80
1:B:602:ARG:HG3	1:B:604:SER:N	1.96	0.80
1:B:414:VAL:HG11	1:B:478:PHE:CE2	2.15	0.80
1:B:373:ARG:CD	4:B:1273:HOH:O	2.29	0.80
1:A:602:ARG:HG2	1:A:603:GLY:N	1.97	0.79
1:B:310:SER:HB3	1:B:592:VAL:CG1	2.13	0.79
1:B:466:GLU:HB2	1:B:469:LYS:HE3	1.64	0.79
1:A:449:CYS:SG	1:A:452:CYS:HB2	2.23	0.79
1:A:330:MET:CE	1:A:783:ALA:HB1	2.12	0.79
1:A:487:GLU:OE2	1:A:607:VAL:HG12	1.83	0.78
1:A:855:LEU:O	1:A:858:THR:HB	1.84	0.78
1:B:546:ARG:NH2	4:B:1259:HOH:O	2.17	0.78
1:B:598:ARG:HE	1:B:600:ASP:HB3	1.48	0.77
1:A:602:ARG:HG3	1:A:604:SER:H	1.48	0.77
1:B:414:VAL:CG1	1:B:478:PHE:CZ	2.68	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:742:ARG:HH11	1:A:742:ARG:CG	1.88	0.76
1:A:723:MET:HG2	1:A:842:PRO:HG3	1.70	0.74
1:B:449:CYS:SG	1:B:452:CYS:HB2	2.29	0.73
1:B:721:ILE:O	1:B:721:ILE:HD12	1.88	0.73
1:B:297:ASP:OD2	1:B:584:ARG:NH2	2.20	0.72
1:A:602:ARG:CG	1:A:603:GLY:H	1.94	0.72
1:B:457:MET:SD	1:B:478:PHE:CE2	2.82	0.72
1:B:666:SER:OG	1:B:803:HIS:HE1	1.72	0.72
1:A:487:GLU:CG	1:A:607:VAL:HG12	2.19	0.72
1:B:373:ARG:HD2	4:B:1273:HOH:O	1.87	0.72
1:A:602:ARG:CG	1:A:603:GLY:N	2.53	0.71
1:B:414:VAL:CG1	1:B:478:PHE:CE2	2.74	0.71
1:B:607:VAL:HG13	1:B:608:VAL:HG13	1.72	0.71
1:A:414:VAL:CG1	1:A:478:PHE:CE1	2.71	0.70
1:A:487:GLU:OE2	1:A:575:TYR:OH	2.08	0.70
1:A:633:GLY:N	1:A:636:HIS:HD2	1.87	0.70
1:B:584:ARG:HD3	1:B:594:ASP:OD2	1.92	0.70
1:A:600:ASP:CG	1:A:601:GLN:N	2.44	0.70
1:B:604:SER:OG	1:B:605:GLY:N	2.24	0.69
1:A:775:ARG:NH2	1:A:844:VAL:O	2.24	0.69
1:B:373:ARG:HD3	4:B:1273:HOH:O	1.92	0.69
1:A:298:PRO:CB	1:A:584:ARG:HH12	2.04	0.68
1:A:607:VAL:HG13	1:A:608:VAL:HG13	1.76	0.67
1:B:721:ILE:HD11	1:B:729:ILE:CG1	2.25	0.67
1:B:600:ASP:CG	1:B:601:GLN:N	2.48	0.67
1:A:292:THR:HG22	1:A:293:THR:HG23	1.75	0.67
1:B:310:SER:HB3	1:B:592:VAL:HG11	1.77	0.66
1:A:298:PRO:HB2	1:A:584:ARG:NH1	2.02	0.66
1:A:416:ALA:HA	1:A:418:GLN:NE2	2.09	0.66
1:B:313:VAL:HG23	1:B:591:THR:HG22	1.78	0.66
1:B:717:HIS:CD2	1:B:719:GLN:HE21	2.14	0.65
1:A:457:MET:CG	1:A:478:PHE:CD1	2.74	0.65
1:A:457:MET:HG3	1:A:478:PHE:CG	2.31	0.65
1:B:633:GLY:N	1:B:636:HIS:HD2	1.92	0.65
1:A:384:TRP:CH2	1:A:555:VAL:HG13	2.32	0.65
1:B:598:ARG:HG2	1:B:600:ASP:HB3	1.78	0.65
1:B:815:LEU:HD21	1:B:837:SER:HA	1.77	0.65
1:B:790:VAL:HG21	1:B:885:VAL:CG2	2.27	0.64
1:A:315:ALA:HB2	1:A:591:THR:HG21	1.80	0.63
1:A:522:ARG:HD2	4:A:1108:HOH:O	1.97	0.63
1:A:406:ASN:ND2	1:A:425:ARG:HG2	2.14	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:603:GLY:O	1:B:604:SER:HB3	1.99	0.63
1:B:456:MET:HB2	1:B:477:TRP:CZ3	2.33	0.63
1:A:457:MET:HG3	1:A:478:PHE:CE1	2.33	0.63
1:A:858:THR:HG22	1:A:861:ARG:H	1.62	0.63
1:B:734:ARG:NH1	2:B:1001:GTP:O2G	2.29	0.62
1:A:414:VAL:CG1	1:A:478:PHE:CZ	2.82	0.62
2:A:1001:GTP:C5'	2:A:1001:GTP:O2B	2.47	0.62
1:B:863:THR:O	1:B:867:ASN:ND2	2.31	0.61
1:A:666:SER:OG	1:A:803:HIS:HE1	1.83	0.61
1:A:474:ARG:NH1	2:A:1001:GTP:H2'	2.16	0.61
2:A:1001:GTP:O2B	2:A:1001:GTP:O1G	2.18	0.61
1:B:752:ASN:ND2	4:B:1330:HOH:O	2.33	0.61
1:A:544:ILE:HG21	1:A:601:GLN:CG	2.31	0.61
1:A:541:ASP:HB3	1:A:603:GLY:HA2	1.83	0.61
1:A:602:ARG:CG	1:A:604:SER:H	2.13	0.61
1:A:416:ALA:HB3	1:A:478:PHE:CD2	2.35	0.60
1:B:766:MET:HG3	1:B:808:TRP:CZ2	2.36	0.60
1:B:800:TRP:HB2	2:B:1001:GTP:O2B	2.02	0.59
1:B:684:HIS:HD2	4:B:1286:HOH:O	1.85	0.59
1:A:297:ASP:OD1	1:A:298:PRO:O	2.20	0.59
1:B:761:LYS:HG2	1:B:794:PRO:HG3	1.84	0.59
1:A:502:ARG:NH2	1:A:663:MET:O	2.30	0.59
1:B:414:VAL:HG11	1:B:478:PHE:CZ	2.35	0.59
1:B:666:SER:OG	1:B:803:HIS:CE1	2.55	0.59
1:A:460:ARG:HB2	1:A:476:ILE:HD11	1.86	0.58
1:A:578:LYS:HZ1	1:A:602:ARG:H	1.49	0.58
1:B:466:GLU:HB2	1:B:469:LYS:CE	2.34	0.57
1:A:506:GLY:HA3	1:A:660:VAL:CG1	2.34	0.57
1:A:576:ARG:O	1:A:599:GLU:HA	2.02	0.57
1:B:717:HIS:HD2	1:B:719:GLN:HE21	1.50	0.57
1:B:522:ARG:HD2	4:B:1113:HOH:O	2.04	0.57
1:A:487:GLU:HG2	1:A:607:VAL:CG1	2.35	0.57
1:A:762:ALA:HB1	1:A:809:MET:HE3	1.85	0.57
1:A:446:ARG:HG2	1:A:446:ARG:O	2.04	0.56
1:B:598:ARG:HE	1:B:600:ASP:CB	2.16	0.56
1:B:416:ALA:HB3	1:B:478:PHE:HE1	1.69	0.56
1:A:598:ARG:HG2	1:A:600:ASP:HB2	1.86	0.56
1:A:633:GLY:N	1:A:636:HIS:CD2	2.69	0.56
1:B:384:TRP:CH2	1:B:555:VAL:HG13	2.41	0.55
1:B:406:ASN:CG	1:B:425:ARG:HH11	2.08	0.55
1:B:312:GLU:OE2	1:B:590:LYS:HE3	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1001:GTP:O2B	2:A:1001:GTP:H5'	2.07	0.54
1:B:721:ILE:HD11	1:B:729:ILE:HD11	1.89	0.54
1:A:710:GLN:NE2	4:A:1120:HOH:O	2.39	0.54
1:A:544:ILE:HG21	1:A:601:GLN:HG3	1.88	0.54
1:B:632:ILE:HD12	1:B:685:PHE:CZ	2.43	0.54
1:A:422:SER:OG	1:A:426:GLU:OE2	2.25	0.54
1:A:752:ASN:HB2	1:A:755:ASP:H	1.74	0.53
1:A:487:GLU:CG	1:A:607:VAL:CG1	2.87	0.53
1:A:617:ASN:O	1:A:621:GLN:HG2	2.09	0.53
1:A:297:ASP:C	1:A:298:PRO:O	2.43	0.52
1:A:330:MET:HE1	1:A:783:ALA:CB	2.28	0.52
1:B:284:GLN:CG	4:B:1297:HOH:O	2.30	0.52
1:A:808:TRP:HD1	1:A:809:MET:HE3	1.75	0.52
1:B:723:MET:HG2	1:B:842:PRO:HG3	1.91	0.52
1:A:544:ILE:HG21	1:A:601:GLN:HG2	1.92	0.51
1:B:546:ARG:NH1	1:B:550:GLU:OE2	2.44	0.51
1:A:707:ASP:HB3	1:A:710:GLN:HE21	1.74	0.51
1:B:773:HIS:H	1:B:773:HIS:CD2	2.28	0.51
1:A:578:LYS:HZ1	1:A:602:ARG:N	2.09	0.51
1:A:296:LYS:O	1:A:297:ASP:HB3	2.11	0.51
1:B:406:ASN:ND2	1:B:425:ARG:HH11	2.09	0.51
1:A:493:PHE:HD2	4:A:1270:HOH:O	1.92	0.51
1:B:315:ALA:HB2	1:B:591:THR:OG1	2.10	0.51
1:B:425:ARG:HH21	1:B:429:ASN:CG	2.14	0.51
1:A:474:ARG:HH12	2:A:1001:GTP:H2'	1.74	0.51
1:B:602:ARG:HD2	1:B:609:THR:HG23	1.93	0.50
1:A:808:TRP:CD1	1:A:809:MET:CE	2.95	0.50
1:B:421:TRP:O	1:B:422:SER:CB	2.59	0.50
1:B:721:ILE:HD11	1:B:729:ILE:HG13	1.93	0.50
1:B:736:GLN:OE1	1:B:776:ASP:HB2	2.12	0.50
1:B:602:ARG:C	1:B:604:SER:H	2.13	0.50
1:A:330:MET:CE	1:A:783:ALA:CB	2.87	0.50
1:B:441:ARG:HD3	4:B:1302:HOH:O	2.12	0.50
1:B:868:ILE:O	1:B:872:ILE:HG13	2.12	0.49
1:A:297:ASP:CG	1:A:298:PRO:O	2.51	0.49
1:A:808:TRP:HD1	1:A:809:MET:CE	2.25	0.49
1:A:761:LYS:HG2	1:A:794:PRO:HG3	1.93	0.49
1:B:732:PRO:O	1:B:773:HIS:CE1	2.65	0.49
1:B:802:ILE:HD12	1:B:802:ILE:C	2.32	0.49
1:A:414:VAL:CG1	1:A:478:PHE:CD1	2.95	0.49
1:A:633:GLY:HA3	1:A:635:GLN:NE2	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:633:GLY:N	1:B:636:HIS:CD2	2.69	0.49
1:A:415:PHE:O	1:A:417:GLU:N	2.44	0.49
1:A:414:VAL:HG11	1:A:478:PHE:CD1	2.46	0.49
1:B:602:ARG:HB2	1:B:602:ARG:HH11	1.77	0.48
1:A:603:GLY:CA	4:A:1280:HOH:O	2.42	0.48
1:B:600:ASP:C	1:B:601:GLN:OE1	2.52	0.48
1:B:561:GLY:O	1:B:565:MET:HG3	2.13	0.48
1:B:632:ILE:HD12	1:B:685:PHE:CE1	2.48	0.48
1:B:861:ARG:HD2	1:B:864:TRP:CZ3	2.48	0.48
1:A:634:PRO:HD2	1:A:635:GLN:NE2	2.20	0.48
1:B:721:ILE:HD13	1:B:842:PRO:CG	2.43	0.47
1:A:415:PHE:CD2	1:A:415:PHE:N	2.70	0.47
1:A:578:LYS:NZ	1:A:602:ARG:H	2.12	0.47
1:B:721:ILE:HD11	1:B:729:ILE:CD1	2.44	0.47
1:A:552:GLU:O	1:A:555:VAL:HG22	2.15	0.47
1:B:721:ILE:HG21	1:B:844:VAL:HG12	1.95	0.47
1:A:734:ARG:NH1	2:A:1001:GTP:O3G	2.43	0.47
1:A:657:GLU:HG3	4:A:1182:HOH:O	2.14	0.47
1:B:373:ARG:HH11	1:B:373:ARG:HB3	1.79	0.47
1:A:443:ASN:OD1	1:A:446:ARG:NH1	2.47	0.47
1:B:278:LYS:HD2	1:B:577:HIS:CD2	2.50	0.47
1:A:742:ARG:CG	1:A:742:ARG:NH1	2.57	0.47
1:B:332:LYS:HB3	1:B:333:PRO:HD3	1.97	0.47
1:B:414:VAL:CG1	1:B:478:PHE:CE1	2.98	0.47
1:B:500:LEU:HD22	1:B:614:THR:CG2	2.44	0.47
1:A:341:THR:OG1	1:B:703:HIS:CD2	2.68	0.47
1:A:487:GLU:HG2	1:A:607:VAL:HG12	1.92	0.47
1:A:789:PRO:HG2	1:A:792:TRP:CE2	2.50	0.47
1:A:832:LYS:O	1:A:834:PRO:HD3	2.15	0.47
1:B:617:ASN:O	1:B:621:GLN:HG2	2.14	0.46
1:A:666:SER:OG	1:A:803:HIS:CE1	2.68	0.46
1:B:322:LEU:HD13	1:B:322:LEU:N	2.31	0.46
1:B:604:SER:HA	1:B:609:THR:OG1	2.16	0.46
1:B:479:MET:HB2	1:B:484:ARG:CD	2.46	0.46
1:B:603:GLY:O	1:B:604:SER:CB	2.63	0.46
1:B:685:PHE:O	1:B:689:MET:HG3	2.15	0.46
1:B:425:ARG:HH21	1:B:429:ASN:CB	2.28	0.46
1:A:742:ARG:NH2	2:A:1001:GTP:O1B	2.48	0.46
1:A:808:TRP:CD1	1:A:809:MET:HE3	2.49	0.46
1:B:598:ARG:NE	1:B:600:ASP:HB3	2.25	0.46
1:A:519:TYR:HB3	1:A:828:TRP:CE2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:416:ALA:HB3	1:B:478:PHE:CE1	2.50	0.46
1:A:874:GLN:O	1:A:878:VAL:HG23	2.16	0.46
1:B:789:PRO:HG2	1:B:792:TRP:CE2	2.50	0.46
1:A:506:GLY:HA3	1:A:660:VAL:HG13	1.99	0.45
1:A:307:TYR:HA	1:A:596:ILE:HG22	1.98	0.45
1:B:456:MET:CB	1:B:477:TRP:CZ3	2.99	0.45
1:B:460:ARG:NH2	1:B:463:LYS:HD2	2.31	0.45
1:B:752:ASN:OD1	1:B:752:ASN:N	2.49	0.45
1:A:601:GLN:HE21	1:A:602:ARG:CZ	2.30	0.45
1:B:721:ILE:CD1	1:B:729:ILE:CG1	2.92	0.45
1:A:809:MET:HA	1:A:809:MET:CE	2.47	0.45
1:A:463:LYS:NZ	2:A:1001:GTP:O3B	2.50	0.45
1:A:446:ARG:CG	1:A:446:ARG:O	2.64	0.45
1:A:601:GLN:HE21	1:A:602:ARG:NH2	2.15	0.45
1:B:721:ILE:HD12	1:B:721:ILE:C	2.36	0.45
1:A:406:ASN:HD22	1:A:425:ARG:HG2	1.79	0.45
1:A:694:LYS:HB2	1:A:694:LYS:HE3	1.78	0.45
1:B:639:GLN:O	1:B:641:PRO:HD3	2.17	0.45
1:A:349:THR:O	1:A:353:GLN:HG3	2.17	0.45
1:A:689:MET:O	1:A:690:SER:HB2	2.16	0.45
1:B:419:ASN:O	1:B:420:GLN:HB3	2.17	0.45
1:A:602:ARG:HD3	1:A:609:THR:HG23	1.99	0.44
1:B:430:ASP:HA	1:B:431:PRO:HD2	1.80	0.44
1:A:773:HIS:CD2	1:A:773:HIS:H	2.36	0.44
1:A:809:MET:CA	1:A:809:MET:HE2	2.47	0.44
1:A:430:ASP:HA	1:A:431:PRO:HD2	1.92	0.44
1:A:304:THR:O	1:A:598:ARG:HD2	2.16	0.44
1:A:578:LYS:NZ	1:A:602:ARG:N	2.66	0.44
1:B:441:ARG:NE	4:B:1302:HOH:O	2.51	0.44
1:A:479:MET:HB2	1:A:484:ARG:CD	2.48	0.44
1:A:546:ARG:O	1:A:550:GLU:HG3	2.18	0.44
1:A:794:PRO:HB2	1:A:809:MET:HE1	1.98	0.44
1:A:512:SER:O	1:A:803:HIS:HD2	2.00	0.44
1:B:297:ASP:OD1	1:B:298:PRO:HD2	2.17	0.44
1:B:636:HIS:HE1	4:B:1307:HOH:O	2.00	0.44
1:A:556:LEU:HA	1:A:556:LEU:HD23	1.88	0.44
1:B:635:GLN:HG2	1:B:636:HIS:CE1	2.52	0.44
1:B:805:LYS:HB2	4:B:1265:HOH:O	2.17	0.44
1:A:322:LEU:HD12	1:A:322:LEU:O	2.17	0.44
1:A:823:ILE:HA	1:A:829:MET:HG2	2.00	0.43
1:B:602:ARG:C	1:B:604:SER:N	2.71	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:528:GLN:O	1:A:662:ARG:NH2	2.35	0.43
1:A:298:PRO:O	1:A:300:HIS:N	2.51	0.43
1:A:418:GLN:HG2	1:A:481:LEU:HD22	2.00	0.43
1:A:278:LYS:HD2	1:A:577:HIS:CD2	2.53	0.43
1:B:493:PHE:HD2	4:B:1127:HOH:O	2.02	0.43
1:A:809:MET:N	1:A:809:MET:HE2	2.34	0.43
1:A:403:LYS:O	1:A:407:SER:OG	2.37	0.43
1:A:584:ARG:HA	1:A:585:PRO:HD3	1.91	0.43
1:A:469:LYS:HE2	1:A:712:PRO:HB2	2.01	0.43
1:B:789:PRO:HD2	1:B:792:TRP:CE3	2.54	0.42
1:A:292:THR:HG22	1:A:293:THR:N	2.34	0.42
1:B:313:VAL:CG2	1:B:591:THR:HG22	2.47	0.42
1:A:773:HIS:HA	1:A:843:TYR:HA	2.01	0.42
1:A:869:TYR:OH	1:A:888:MET:HB2	2.19	0.42
1:A:800:TRP:N	2:A:1001:GTP:O1B	2.38	0.42
1:B:746:SER:HA	1:B:747:PRO:HD3	1.91	0.42
1:A:656:GLY:O	1:A:660:VAL:HB	2.20	0.42
1:B:339:ASN:HD22	1:B:339:ASN:HA	1.67	0.42
1:B:541:ASP:HB3	1:B:603:GLY:HA2	2.02	0.42
1:A:479:MET:HB2	1:A:484:ARG:NE	2.35	0.42
1:A:437:VAL:HG11	1:A:485:TYR:CD2	2.54	0.42
1:A:772:PHE:O	1:A:778:ARG:HD3	2.20	0.42
1:B:721:ILE:CD1	1:B:729:ILE:HG13	2.50	0.42
1:B:414:VAL:HG13	1:B:478:PHE:CZ	2.52	0.41
1:A:789:PRO:HG2	1:A:792:TRP:CZ2	2.55	0.41
1:A:786:SER:HA	1:A:885:VAL:O	2.20	0.41
1:A:297:ASP:O	1:A:298:PRO:C	2.53	0.41
1:A:442:GLU:HG3	4:A:1248:HOH:O	2.20	0.41
1:A:601:GLN:NE2	1:A:602:ARG:CZ	2.84	0.41
1:B:415:PHE:O	1:B:417:GLU:N	2.52	0.41
1:B:601:GLN:HG2	1:B:601:GLN:O	2.21	0.41
1:B:322:LEU:HB2	1:B:746:SER:O	2.20	0.41
1:B:723:MET:SD	1:B:835:ILE:HD11	2.60	0.41
1:B:766:MET:HG3	1:B:808:TRP:CD2	2.55	0.41
1:A:477:TRP:HB2	1:A:605:GLY:HA3	2.03	0.41
1:A:457:MET:CG	1:A:478:PHE:CG	3.02	0.40
1:A:350:PRO:HG2	1:A:591:THR:HG21	2.04	0.40
1:B:666:SER:O	1:B:669:ASP:HB2	2.21	0.40
1:B:833:THR:HA	1:B:834:PRO:HD3	1.93	0.40
1:B:409:ALA:HB1	1:B:410:ALA:H	1.51	0.40
1:B:602:ARG:NH1	1:B:602:ARG:HB2	2.35	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:809:MET:HA	1:A:809:MET:HE2	2.04	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	608/639 (95%)	572 (94%)	28 (5%)	8 (1%)	12	15
1	B	608/639 (95%)	570 (94%)	28 (5%)	10 (2%)	9	11
All	All	1216/1278 (95%)	1142 (94%)	56 (5%)	18 (2%)	10	12

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	297	ASP
1	A	416	ALA
1	B	409	ALA
1	B	416	ALA
1	B	420	GLN
1	B	752	ASN
1	A	299	GLU
1	A	587	ALA
1	A	604	SER
1	A	752	ASN
1	B	422	SER
1	B	604	SER
1	B	408	ASN
1	A	602	ARG
1	A	413	ALA
1	B	601	GLN
1	B	413	ALA

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Mol	Chain	Res	Type
1	B	603	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	516/544 (95%)	494 (96%)	22 (4%)	29	43
1	B	514/544 (94%)	494 (96%)	20 (4%)	32	48
All	All	1030/1088 (95%)	988 (96%)	42 (4%)	30	45

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

Mol	Chain	Res	Type
1	A	285	LYS
1	A	286	LEU
1	A	348	THR
1	A	363	THR
1	A	407	SER
1	A	411	LEU
1	A	415	PHE
1	A	418	GLN
1	A	459	LYS
1	A	466	GLU
1	A	555	VAL
1	A	635	GLN
1	A	660	VAL
1	A	724	LYS
1	A	742	ARG
1	A	778	ARG
1	A	811	THR
1	A	839	THR
1	A	847	ARG
1	A	858	THR
1	A	859	ARG
1	A	861	ARG

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Mol	Chain	Res	Type
1	B	292	THR
1	B	322	LEU
1	B	328	LYS
1	B	355	ARG
1	B	415	PHE
1	B	418	GLN
1	B	441	ARG
1	B	457	MET
1	B	469	LYS
1	B	555	VAL
1	B	602	ARG
1	B	659	ARG
1	B	669	ASP
1	B	724	LYS
1	B	733	CYS
1	B	752	ASN
1	B	778	ARG
1	B	847	ARG
1	B	861	ARG
1	B	867	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	406	ASN
1	A	408	ASN
1	A	418	GLN
1	A	419	ASN
1	A	450	HIS
1	A	551	ASN
1	A	635	GLN
1	A	636	HIS
1	A	687	ASN
1	A	703	HIS
1	A	710	GLN
1	A	717	HIS
1	A	803	HIS
1	B	339	ASN
1	B	495	ASN
1	B	551	ASN
1	B	563	HIS
1	B	636	HIS

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Mol	Chain	Res	Type
1	B	684	HIS
1	B	687	ASN
1	B	703	HIS
1	B	717	HIS
1	B	765	GLN
1	B	803	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are 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	$\# Z > 2$
2	GTP	B	1001	-	26,34,34	1.89	6 (23%)	33,54,54	1.96	7 (21%)
2	GTP	A	1001	-	26,34,34	2.08	6 (23%)	33,54,54	2.36	8 (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
2	GTP	B	1001	-	-	6/18/38/38	0/3/3/3
2	GTP	A	1001	-	-	8/18/38/38	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	GTP	O4'-C1'	6.97	1.50	1.41
2	B	1001	GTP	O4'-C1'	5.15	1.48	1.41
2	A	1001	GTP	C6-N1	4.34	1.40	1.33
2	B	1001	GTP	C6-N1	3.79	1.39	1.33
2	A	1001	GTP	C2-N1	3.31	1.41	1.35
2	B	1001	GTP	C2-N1	3.12	1.40	1.35
2	B	1001	GTP	PG-O3G	2.76	1.65	1.54
2	A	1001	GTP	PG-O3G	2.61	1.64	1.54
2	B	1001	GTP	PA-O5'	2.21	1.68	1.59
2	B	1001	GTP	C5'-C4'	2.16	1.58	1.51
2	A	1001	GTP	O2'-C2'	2.11	1.47	1.43
2	A	1001	GTP	PG-O2G	2.09	1.62	1.54

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	GTP	N3-C2-N1	-6.74	118.24	127.22
2	B	1001	GTP	N3-C2-N1	-5.83	119.44	127.22
2	A	1001	GTP	PB-O3B-PG	-5.78	112.98	132.83
2	A	1001	GTP	C2-N3-C4	5.58	121.73	115.36
2	B	1001	GTP	C2-N3-C4	5.44	121.57	115.36
2	A	1001	GTP	PA-O3A-PB	-4.23	118.31	132.83
2	B	1001	GTP	PB-O3B-PG	-3.91	119.41	132.83
2	A	1001	GTP	C6-N1-C2	3.24	121.07	115.93
2	A	1001	GTP	C5-C6-N1	-3.13	119.16	123.43
2	A	1001	GTP	C6-C5-C4	-2.67	118.25	120.80
2	B	1001	GTP	C6-C5-C4	-2.55	118.36	120.80
2	B	1001	GTP	N2-C2-N3	2.41	121.72	117.79
2	A	1001	GTP	N2-C2-N3	2.35	121.62	117.79
2	B	1001	GTP	C5-C6-N1	-2.13	120.51	123.43
2	B	1001	GTP	C6-N1-C2	2.11	119.28	115.93

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	GTP	C5'-O5'-PA-O1A
2	A	1001	GTP	C5'-O5'-PA-O2A
2	B	1001	GTP	C5'-O5'-PA-O3A
2	B	1001	GTP	C4'-C5'-O5'-PA
2	A	1001	GTP	C4'-C5'-O5'-PA
2	B	1001	GTP	PB-O3A-PA-O1A
2	A	1001	GTP	PB-O3A-PA-O5'
2	B	1001	GTP	PB-O3A-PA-O5'
2	A	1001	GTP	PA-O3A-PB-O2B
2	B	1001	GTP	C5'-O5'-PA-O1A
2	A	1001	GTP	C3'-C4'-C5'-O5'
2	B	1001	GTP	PG-O3B-PB-O2B
2	A	1001	GTP	O4'-C4'-C5'-O5'
2	A	1001	GTP	C5'-O5'-PA-O3A

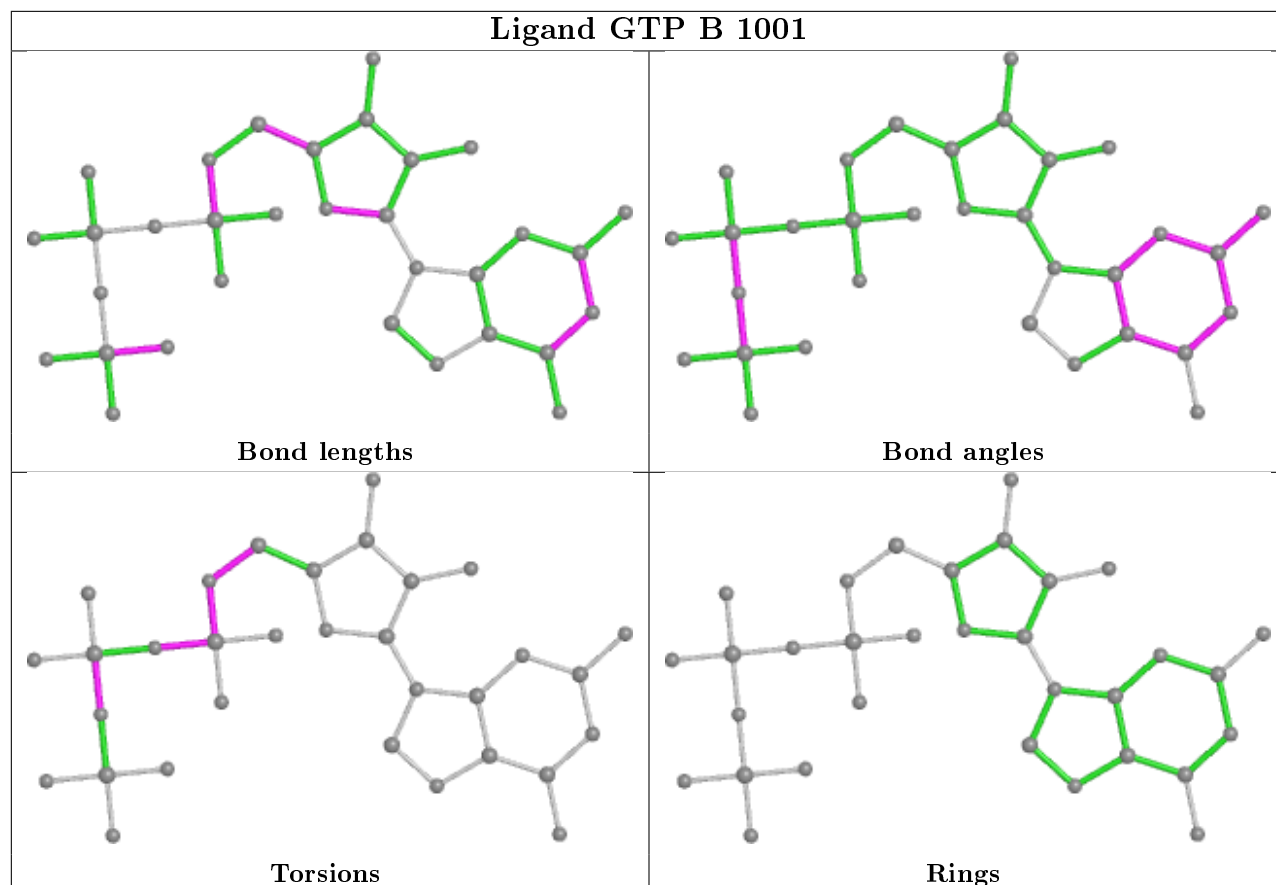
There are no ring outliers.

2 monomers are involved in 11 short contacts:

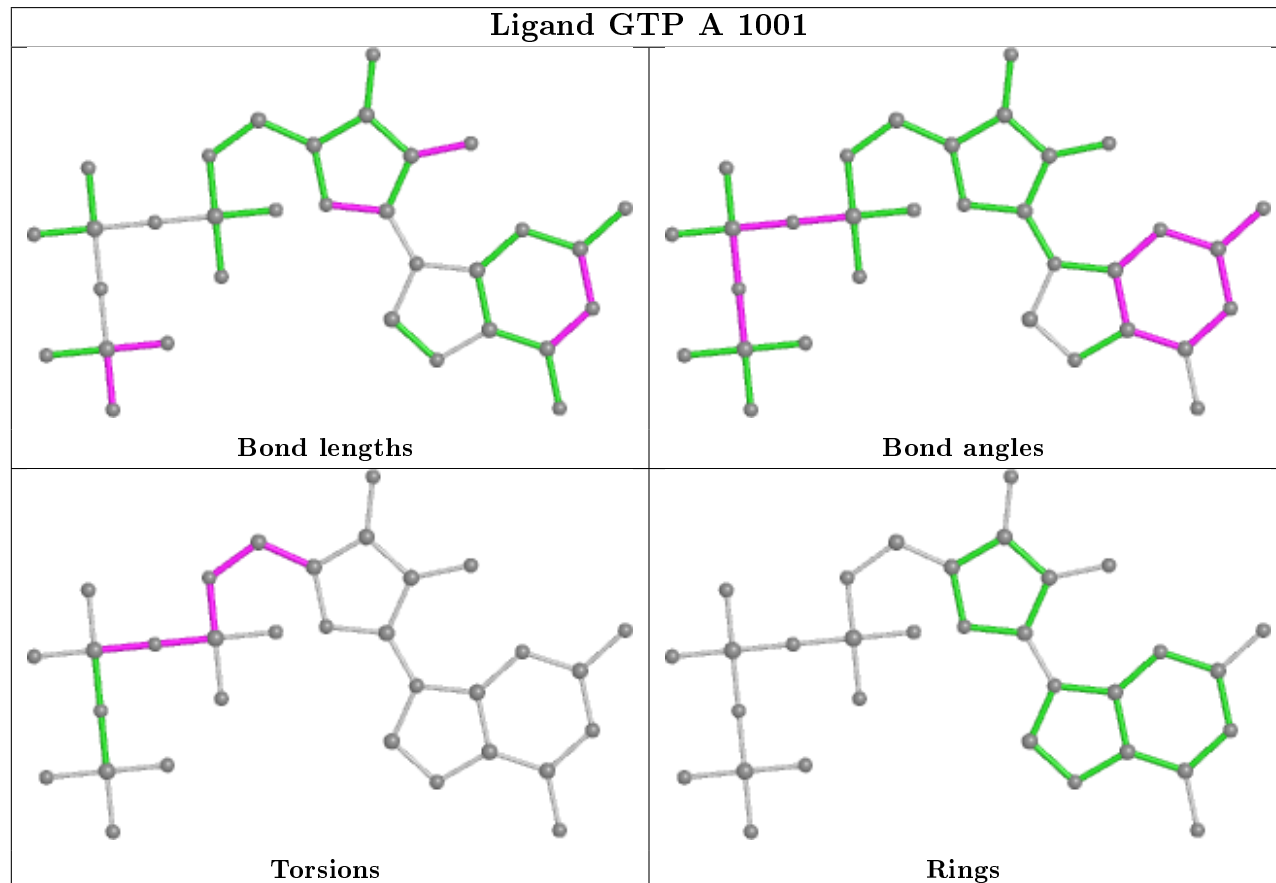
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1001	GTP	2	0
2	A	1001	GTP	9	0

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.

Ligand GTP B 1001



Ligand GTP A 1001



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, 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	612/639 (95%)	0.05	18 (2%) 51 53	15, 33, 60, 81	0
1	B	612/639 (95%)	0.02	15 (2%) 57 59	16, 33, 57, 76	0
All	All	1224/1278 (95%)	0.04	33 (2%) 54 56	15, 33, 58, 81	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	416	ALA	9.0
1	A	415	PHE	6.1
1	B	416	ALA	5.2
1	A	412	GLY	4.6
1	A	413	ALA	4.4
1	B	751	TRP	4.3
1	A	420	GLN	4.2
1	B	411	LEU	4.1
1	A	750	GLY	3.9
1	A	419	ASN	3.8
1	B	415	PHE	3.7
1	B	414	VAL	3.7
1	A	275	ASN	3.1
1	A	478	PHE	2.7
1	A	751	TRP	2.7
1	A	411	LEU	2.6
1	A	889	THR	2.6
1	B	750	GLY	2.6
1	B	413	ALA	2.5
1	A	602	ARG	2.5
1	B	425	ARG	2.5
1	B	410	ALA	2.4
1	A	603	GLY	2.4
1	B	422	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	276	GLN	2.3
1	A	408	ASN	2.3
1	B	419	ASN	2.2
1	A	418	GLN	2.2
1	B	420	GLN	2.2
1	B	315	ALA	2.1
1	A	587	ALA	2.1
1	A	588	GLU	2.0
1	B	292	THR	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 monosaccharides in this entry.

6.4 Ligands [i](#)

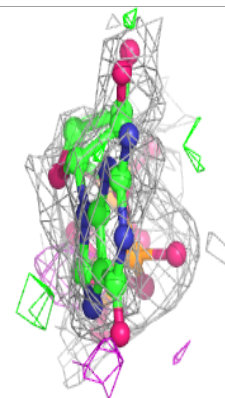
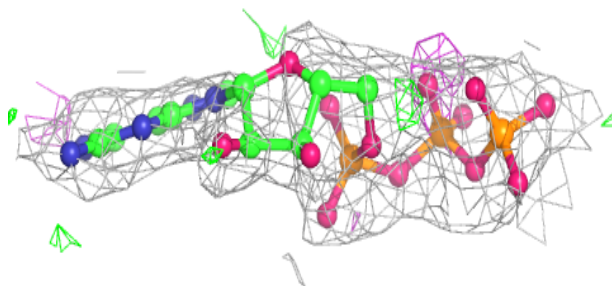
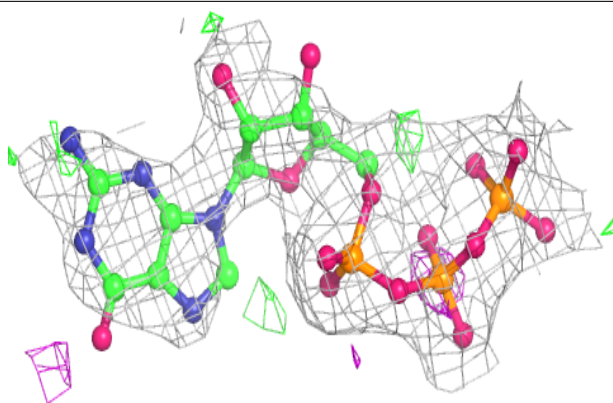
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	GTP	A	1001	32/32	0.90	0.19	44,72,78,80	0
2	GTP	B	1001	32/32	0.90	0.20	52,71,78,78	0
3	ZN	A	1003	1/1	0.98	0.09	36,36,36,36	1
3	ZN	A	1002	1/1	0.99	0.08	33,33,33,33	0
3	ZN	B	1003	1/1	0.99	0.07	37,37,37,37	1
3	ZN	B	1002	1/1	0.99	0.07	31,31,31,31	0

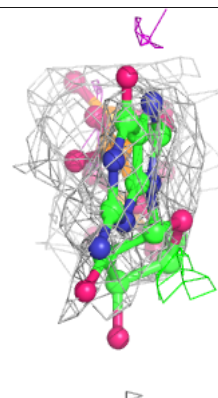
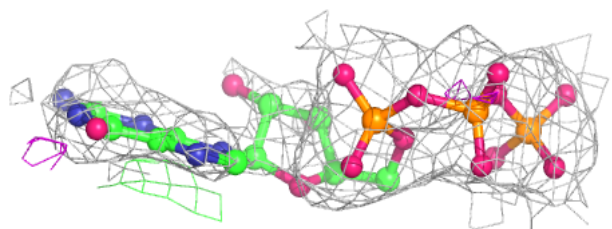
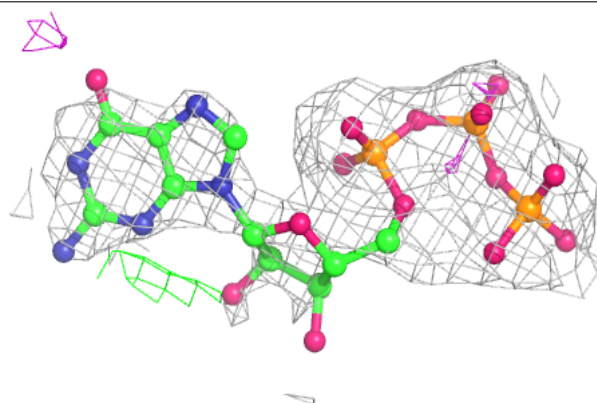
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around GTP A 1001:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around GTP B 1001:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.