



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 06:47 PM UTC

PDB ID : 1S49 / pdb_00001s49
Title : Crystal Structure of RNA-dependent RNA polymerase construct 1 (residues 71-679) from bovine viral diarrhea virus complexed with GTP
Authors : Choi, K.H.; Groarke, J.M.; Young, D.C.; Kuhn, R.J.; Smith, J.L.; Pevear, D.C.; Rossmann, M.G.
Deposited on : 2004-01-15
Resolution : 3.00 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

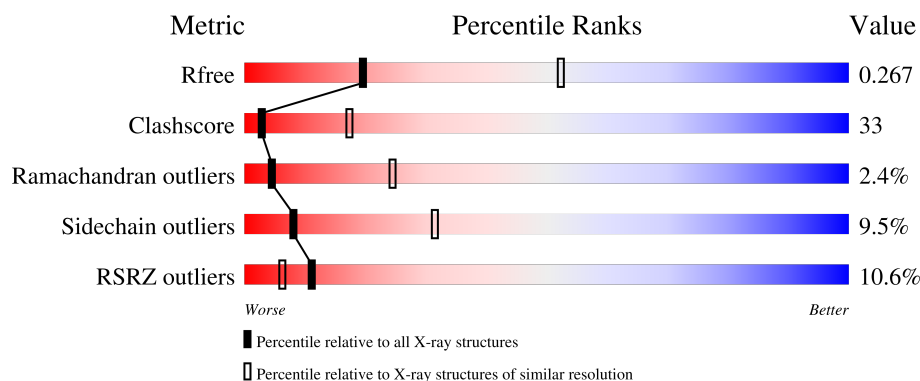
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	609	10% 46% 41% 9% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4772 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.

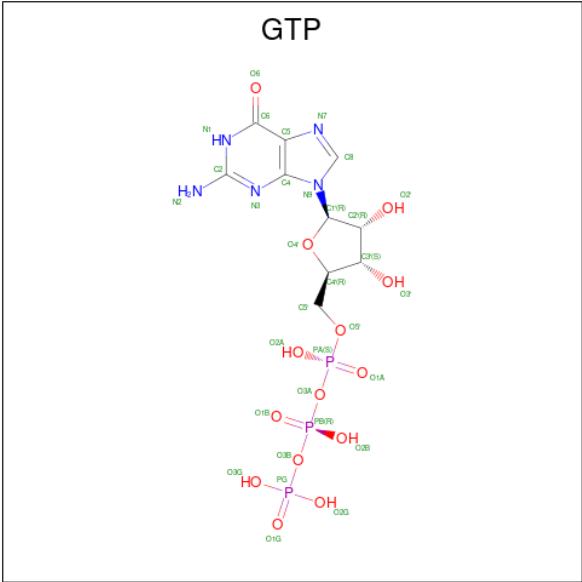
- Molecule 1 is a protein called RNA-dependent RNA polymerase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	588	4740	3011	837	875	5	12	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	115	MSE	MET	modified residue	UNP P19711
A	143	MSE	MET	modified residue	UNP P19711
A	302	MSE	MET	modified residue	UNP P19711
A	384	MSE	MET	modified residue	UNP P19711
A	416	MSE	MET	modified residue	UNP P19711
A	422	MSE	MET	modified residue	UNP P19711
A	423	MSE	MET	modified residue	UNP P19711
A	467	MSE	MET	modified residue	UNP P19711
A	485	MSE	MET	modified residue	UNP P19711
A	514	MSE	MET	modified residue	UNP P19711
A	526	MSE	MET	modified residue	UNP P19711
A	552	MSE	MET	modified residue	UNP P19711

- Molecule 2 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).



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

4 Data and refinement statistics

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, α , β , γ	205.73Å 205.73Å 99.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.78 – 3.00 29.78 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.6 (29.78-3.00) 98.4 (29.78-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 2.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.264 , 0.266 0.263 , 0.267	Depositor DCC
R_{free} test set	5635 reflections (7.79%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.106	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	4772	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.66% 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	0.52	1/4830 (0.0%)	1.20	44/6497 (0.7%)

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	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	531	ASP	CA-C	-6.02	1.45	1.52

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	533	SER	N-CA-C	-28.99	82.24	112.97
1	A	539	THR	N-CA-C	-13.21	92.33	113.19
1	A	673	HIS	N-CA-C	12.15	128.46	110.48
1	A	291	GLU	N-CA-C	11.80	128.29	112.25
1	A	402	GLN	N-CA-C	-9.12	91.37	110.80
1	A	181	ASN	CA-C-N	8.91	130.98	119.84
1	A	181	ASN	C-N-CA	8.91	130.98	119.84
1	A	209	TRP	N-CA-C	-8.52	101.99	112.38
1	A	293	LYS	N-CA-C	-8.08	93.58	110.80
1	A	530	LEU	CA-C-N	-8.00	110.06	121.50
1	A	530	LEU	C-N-CA	-8.00	110.06	121.50
1	A	289	TYR	N-CA-C	7.84	120.93	108.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	387	VAL	N-CA-C	7.67	116.25	107.89
1	A	664	ASP	N-CA-C	-7.64	97.78	109.85
1	A	309	GLN	CA-C-N	-7.60	110.34	119.84
1	A	309	GLN	C-N-CA	-7.60	110.34	119.84
1	A	338	GLU	CA-C-N	7.57	127.63	119.90
1	A	338	GLU	C-N-CA	7.57	127.63	119.90
1	A	540	ALA	N-CA-C	-7.28	104.43	113.38
1	A	212	LEU	N-CA-C	-7.14	103.50	111.28
1	A	531	ASP	N-CA-C	6.96	120.82	109.76
1	A	531	ASP	O-C-N	6.90	131.12	123.04
1	A	318	GLY	N-CA-C	6.83	121.16	112.83
1	A	275	ALA	N-CA-C	-6.70	100.31	109.95
1	A	146	ALA	N-CA-C	-6.25	102.47	110.53
1	A	461	LYS	N-CA-C	-6.16	104.64	111.36
1	A	127	ARG	N-CA-C	-6.07	105.13	112.54
1	A	294	THR	N-CA-C	-6.03	104.82	111.82
1	A	289	TYR	CA-C-N	6.03	127.38	119.84
1	A	289	TYR	C-N-CA	6.03	127.38	119.84
1	A	531	ASP	N-CA-CB	5.96	119.19	109.95
1	A	216	VAL	N-CA-C	5.93	117.99	109.51
1	A	304	ASN	N-CA-C	5.83	117.63	111.28
1	A	232	GLU	N-CA-C	-5.72	105.12	111.36
1	A	376	PHE	N-CA-C	-5.68	104.99	111.07
1	A	553	TYR	N-CA-C	5.59	119.29	111.30
1	A	589	GLY	N-CA-C	-5.52	105.81	112.49
1	A	541	TYR	N-CA-C	-5.52	106.59	113.55
1	A	323	PHE	N-CA-C	5.42	119.44	113.21
1	A	388	PRO	N-CA-C	-5.27	101.97	110.55
1	A	167	GLU	N-CA-C	-5.19	105.74	111.71
1	A	662	ILE	CA-C-N	5.07	125.35	119.92
1	A	662	ILE	C-N-CA	5.07	125.35	119.92
1	A	431	GLY	N-CA-C	-5.05	108.35	114.92

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	531	ASP	Mainchain

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	4740	0	4781	312	1
2	A	32	0	12	3	0
All	All	4772	0	4793	312	1

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 33.

All (312) 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:532:SER:C	1:A:534:GLY:H	1.29	1.18
1:A:538:THR:HB	1:A:540:ALA:H	1.16	1.08
1:A:178:ASN:HD21	1:A:352:GLN:HG3	1.17	1.07
1:A:336:PHE:HB2	1:A:339:PRO:HG3	1.33	1.06
1:A:532:SER:C	1:A:534:GLY:N	2.07	1.06
1:A:106:PHE:HB3	1:A:151:GLU:HG2	1.36	1.06
1:A:380:ILE:HG12	1:A:384:MSE:HE2	1.35	1.04
1:A:678:GLN:HG2	1:A:678:GLN:O	1.56	1.03
1:A:227:LYS:H	1:A:227:LYS:HD2	1.25	1.01
1:A:175:LYS:HD3	1:A:351:THR:HG23	1.43	0.98
1:A:322:LEU:HG	2:A:2030:GTP:C2	2.01	0.94
1:A:283:ARG:H	1:A:283:ARG:CD	1.78	0.93
1:A:313:ILE:HG12	1:A:416:MSE:HE2	1.50	0.91
1:A:227:LYS:H	1:A:227:LYS:CD	1.82	0.87
1:A:229:ASN:C	1:A:229:ASN:HD22	1.82	0.87
1:A:536:ARG:O	1:A:536:ARG:HD3	1.74	0.87
1:A:229:ASN:HD21	1:A:232:GLU:HG3	1.39	0.86
1:A:536:ARG:O	1:A:536:ARG:CD	2.26	0.84
1:A:227:LYS:HD3	1:A:228:LYS:H	1.40	0.84
1:A:381:THR:HA	1:A:384:MSE:HE3	1.59	0.83
1:A:178:ASN:ND2	1:A:352:GLN:HG3	1.93	0.83
1:A:237:GLU:HB3	1:A:240:LEU:HD12	1.60	0.83
1:A:494:ILE:O	1:A:500:THR:HG22	1.78	0.82
1:A:277:ASP:OD1	1:A:279:VAL:HG12	1.83	0.79
1:A:243:GLN:HG3	1:A:247:ASP:OD2	1.83	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:531:ASP:HB3	1:A:541:TYR:CE1	2.17	0.79
1:A:492:GLU:HA	1:A:501:PRO:HG2	1.64	0.77
1:A:283:ARG:H	1:A:283:ARG:HD3	1.49	0.76
1:A:540:ALA:O	1:A:576:SER:N	2.18	0.76
1:A:349:TRP:CZ3	1:A:475:LYS:HD2	2.20	0.75
1:A:229:ASN:ND2	1:A:232:GLU:HG3	2.02	0.75
1:A:538:THR:HB	1:A:540:ALA:N	1.99	0.74
1:A:254:ILE:HG23	1:A:383:HIS:NE2	2.02	0.74
1:A:536:ARG:HA	1:A:536:ARG:HH11	1.52	0.74
1:A:661:TYR:O	1:A:663:PRO:HD3	1.89	0.72
1:A:218:ARG:C	1:A:220:GLY:H	1.98	0.72
1:A:354:THR:HG22	1:A:356:LYS:H	1.56	0.70
1:A:283:ARG:H	1:A:283:ARG:HD2	1.56	0.70
1:A:292:ALA:HB2	1:A:295:ARG:NH2	2.06	0.70
1:A:209:TRP:O	1:A:213:GLU:HG3	1.91	0.69
1:A:354:THR:HG22	1:A:355:SER:N	2.06	0.69
1:A:531:ASP:OD2	1:A:533:SER:N	2.25	0.69
1:A:419:VAL:HG12	1:A:443:ILE:HD13	1.75	0.69
1:A:152:LYS:O	1:A:154:PRO:HD3	1.93	0.68
1:A:648:LYS:HA	1:A:648:LYS:HZ3	1.59	0.68
1:A:648:LYS:HD3	1:A:659:HIS:HB2	1.75	0.67
1:A:165:PHE:CD1	1:A:262:PRO:HB3	2.29	0.67
1:A:556:ASN:HD22	1:A:559:VAL:H	1.42	0.67
1:A:648:LYS:N	1:A:649:PRO:CD	2.58	0.67
1:A:209:TRP:CD2	1:A:238:LYS:HE3	2.29	0.67
1:A:173:ILE:HD11	1:A:286:VAL:HG21	1.77	0.67
1:A:316:TYR:CZ	1:A:318:GLY:HA3	2.30	0.66
1:A:639:GLY:HA3	1:A:648:LYS:HE2	1.78	0.66
1:A:101:LEU:HA	1:A:104:ILE:HD13	1.78	0.65
1:A:648:LYS:N	1:A:649:PRO:HD3	2.11	0.65
1:A:532:SER:CA	1:A:534:GLY:H	2.07	0.65
1:A:404:GLY:O	1:A:407:GLN:HG3	1.96	0.65
1:A:541:TYR:HA	1:A:576:SER:HB3	1.78	0.65
1:A:291:GLU:CG	1:A:291:GLU:O	2.45	0.65
1:A:531:ASP:HB3	1:A:541:TYR:CZ	2.32	0.64
1:A:226:GLU:OE2	1:A:293:LYS:HG3	1.97	0.64
1:A:138:GLN:OE1	1:A:138:GLN:HA	1.97	0.64
1:A:354:THR:CG2	1:A:355:SER:N	2.61	0.64
1:A:230:ILE:HG12	1:A:296:LEU:HD11	1.79	0.63
1:A:220:GLY:HA3	1:A:230:ILE:HD12	1.79	0.63
1:A:154:PRO:CG	1:A:620:LEU:HD13	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:MSE:HA	1:A:369:TYR:OH	1.97	0.63
1:A:222:ALA:O	1:A:226:GLU:OE1	2.17	0.62
1:A:291:GLU:O	1:A:291:GLU:HG2	1.97	0.62
1:A:164:THR:HA	1:A:167:GLU:HG3	1.80	0.62
1:A:227:LYS:HD3	1:A:228:LYS:N	2.13	0.62
1:A:322:LEU:HG	2:A:2030:GTP:N2	2.14	0.61
1:A:480:THR:CG2	1:A:482:GLY:H	2.13	0.61
1:A:143:MSE:O	1:A:148:ILE:HB	1.99	0.61
1:A:648:LYS:HB3	1:A:648:LYS:HZ2	1.65	0.61
1:A:212:LEU:HD13	1:A:300:LYS:HA	1.83	0.61
1:A:138:GLN:HE21	1:A:539:THR:CG2	2.12	0.60
1:A:164:THR:HG22	1:A:167:GLU:OE2	2.02	0.60
1:A:592:LYS:HB2	1:A:598:ASN:ND2	2.16	0.60
1:A:183:GLU:HB2	1:A:187:LYS:HE2	1.85	0.59
1:A:227:LYS:CD	1:A:227:LYS:N	2.55	0.59
1:A:298:ILE:HG12	1:A:380:ILE:HD12	1.83	0.59
1:A:146:ALA:HB1	1:A:627:THR:HA	1.83	0.59
1:A:303:TYR:O	1:A:307:LYS:HB2	2.02	0.59
1:A:138:GLN:HG3	1:A:539:THR:HG22	1.85	0.58
1:A:336:PHE:CB	1:A:339:PRO:HG3	2.22	0.58
1:A:354:THR:HB	1:A:357:ASP:OD1	2.03	0.58
1:A:173:ILE:CD1	1:A:286:VAL:HG21	2.33	0.58
1:A:674:TYR:CD2	1:A:674:TYR:N	2.70	0.58
1:A:261:ILE:HG22	1:A:392:ALA:HB2	1.86	0.57
1:A:226:GLU:OE1	1:A:291:GLU:O	2.22	0.57
1:A:352:GLN:HG2	1:A:476:PRO:HD2	1.85	0.57
1:A:251:GLY:HA3	1:A:375:LYS:HE2	1.87	0.57
1:A:192:PHE:CE1	1:A:364:ILE:HG23	2.40	0.56
1:A:322:LEU:O	1:A:325:ILE:HG22	2.06	0.56
1:A:146:ALA:CB	1:A:627:THR:HG23	2.36	0.56
1:A:535:GLU:C	1:A:537:GLY:H	2.12	0.56
1:A:160:THR:O	1:A:160:THR:HG22	2.06	0.55
1:A:163:LYS:O	1:A:167:GLU:HG3	2.06	0.55
1:A:221:ALA:O	1:A:295:ARG:NH2	2.39	0.55
1:A:227:LYS:HD2	1:A:227:LYS:N	2.08	0.55
1:A:201:LYS:HG2	1:A:202:HIS:CD2	2.41	0.55
1:A:648:LYS:NZ	1:A:648:LYS:CB	2.70	0.55
1:A:309:GLN:HB2	1:A:310:PRO:CD	2.36	0.55
1:A:640:LYS:HB2	1:A:640:LYS:NZ	2.21	0.55
1:A:218:ARG:C	1:A:220:GLY:N	2.65	0.55
1:A:432:VAL:HG21	1:A:440:VAL:HG21	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:354:THR:CG2	1:A:355:SER:H	2.20	0.54
1:A:92:VAL:HG23	1:A:92:VAL:O	2.07	0.54
1:A:166:HIS:HB3	1:A:170:ARG:HH12	1.72	0.54
1:A:208:THR:C	1:A:210:GLU:N	2.62	0.54
1:A:586:ASP:HB3	1:A:664:ASP:O	2.08	0.54
1:A:146:ALA:HB1	1:A:627:THR:HG23	1.90	0.54
1:A:309:GLN:HB2	1:A:310:PRO:HD3	1.90	0.54
1:A:536:ARG:HD3	1:A:536:ARG:C	2.26	0.54
1:A:322:LEU:HG	2:A:2030:GTP:N3	2.23	0.53
1:A:330:ARG:HD2	1:A:330:ARG:O	2.07	0.53
1:A:670:GLN:HE22	1:A:677:LEU:N	2.06	0.53
1:A:648:LYS:HE3	1:A:657:THR:OG1	2.09	0.53
1:A:165:PHE:O	1:A:168:ALA:HB3	2.09	0.53
1:A:639:GLY:HA2	1:A:645:TRP:CD1	2.43	0.53
1:A:224:PHE:HB2	1:A:259:THR:HG23	1.90	0.52
1:A:519:THR:HG21	1:A:617:LEU:HD11	1.91	0.52
1:A:640:LYS:HB2	1:A:640:LYS:HZ2	1.74	0.52
1:A:309:GLN:O	1:A:311:VAL:N	2.42	0.52
1:A:209:TRP:CG	1:A:238:LYS:HE3	2.44	0.52
1:A:502:VAL:HG23	1:A:516:GLY:HA3	1.92	0.52
1:A:187:LYS:O	1:A:191:ILE:HG13	2.09	0.52
1:A:626:HIS:O	1:A:630:ARG:HG3	2.09	0.52
1:A:229:ASN:C	1:A:229:ASN:ND2	2.52	0.52
1:A:320:THR:HG22	1:A:321:PRO:O	2.10	0.52
1:A:277:ASP:CG	1:A:279:VAL:HG12	2.34	0.52
1:A:506:TRP:CD2	1:A:597:ARG:HD3	2.45	0.52
1:A:322:LEU:HD23	1:A:325:ILE:HG21	1.92	0.51
1:A:337:ASN:C	1:A:339:PRO:HD3	2.36	0.51
1:A:355:SER:HB3	1:A:385:THR:HG23	1.91	0.51
1:A:528:THR:O	1:A:530:LEU:HD12	2.11	0.51
1:A:610:LEU:HD22	1:A:614:ASN:ND2	2.26	0.51
1:A:332:GLU:OE2	1:A:442:ARG:HD2	2.10	0.51
1:A:648:LYS:NZ	1:A:653:ILE:HG22	2.25	0.51
1:A:230:ILE:HD11	1:A:295:ARG:NH1	2.26	0.51
1:A:389:VAL:HG11	1:A:397:TYR:CZ	2.45	0.51
1:A:256:TYR:OH	1:A:290:PRO:HG3	2.11	0.50
1:A:497:CYS:O	1:A:499:HIS:HD2	1.94	0.50
1:A:352:GLN:HG2	1:A:476:PRO:CD	2.41	0.50
1:A:354:THR:HG22	1:A:356:LYS:N	2.26	0.50
1:A:212:LEU:CD1	1:A:299:THR:HG22	2.42	0.50
1:A:292:ALA:H	1:A:295:ARG:HE	1.59	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:307:LYS:O	1:A:308:GLN:C	2.54	0.50
1:A:361:ILE:HD11	1:A:417:LEU:HB2	1.93	0.50
1:A:536:ARG:O	1:A:536:ARG:CG	2.59	0.50
1:A:605:THR:O	1:A:606:GLY:C	2.55	0.50
1:A:322:LEU:HB3	1:A:552:MSE:SE	2.61	0.50
1:A:415:SER:O	1:A:419:VAL:HG22	2.12	0.49
1:A:531:ASP:CG	1:A:532:SER:N	2.67	0.49
1:A:668:THR:O	1:A:669:LEU:HD23	2.12	0.49
1:A:95:GLU:C	1:A:97:ASN:H	2.19	0.49
1:A:506:TRP:CE2	1:A:597:ARG:HD2	2.47	0.49
1:A:227:LYS:HE2	1:A:293:LYS:NZ	2.27	0.49
1:A:301:VAL:HG22	1:A:302:MSE:HG3	1.92	0.49
1:A:138:GLN:HG2	1:A:542:GLU:OE1	2.13	0.49
1:A:173:ILE:HD11	1:A:286:VAL:CG2	2.41	0.49
1:A:224:PHE:HD2	1:A:225:LEU:HD23	1.78	0.49
1:A:322:LEU:HD13	1:A:552:MSE:HE1	1.94	0.49
1:A:292:ALA:HB2	1:A:295:ARG:HH21	1.78	0.49
1:A:438:ASN:OD1	1:A:442:ARG:NH1	2.46	0.49
1:A:338:GLU:N	1:A:339:PRO:HD3	2.28	0.49
1:A:159:GLN:HG2	1:A:281:GLU:CD	2.37	0.49
1:A:309:GLN:O	1:A:310:PRO:C	2.52	0.49
1:A:556:ASN:ND2	1:A:559:VAL:H	2.10	0.49
1:A:648:LYS:HZ3	1:A:648:LYS:CA	2.25	0.48
1:A:126:ASP:HB2	1:A:130:ARG:HB2	1.94	0.48
1:A:648:LYS:HB3	1:A:648:LYS:NZ	2.28	0.48
1:A:452:LEU:HD23	1:A:452:LEU:O	2.14	0.48
1:A:218:ARG:O	1:A:220:GLY:N	2.46	0.48
1:A:531:ASP:OD2	1:A:532:SER:N	2.46	0.48
1:A:541:TYR:HA	1:A:576:SER:CB	2.43	0.48
1:A:636:VAL:O	1:A:637:ALA:C	2.57	0.48
1:A:204:TYR:HB2	1:A:368:TYR:O	2.14	0.48
1:A:112:THR:HG21	1:A:115:MSE:HE3	1.96	0.48
1:A:347:LYS:O	1:A:352:GLN:NE2	2.47	0.48
1:A:263:LYS:HD3	1:A:265:GLU:CD	2.39	0.48
1:A:380:ILE:HG23	1:A:381:THR:N	2.28	0.48
1:A:341:ALA:HB3	1:A:488:ALA:HB3	1.96	0.47
1:A:480:THR:HG23	1:A:482:GLY:H	1.78	0.47
1:A:162:THR:HA	1:A:165:PHE:HB3	1.95	0.47
1:A:528:THR:C	1:A:530:LEU:HD12	2.38	0.47
1:A:207:VAL:HG22	1:A:300:LYS:O	2.14	0.47
1:A:418:ASN:HD22	1:A:445:VAL:HG22	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:531:ASP:OD2	1:A:532:SER:C	2.58	0.47
1:A:581:TYR:OH	1:A:677:LEU:CD1	2.63	0.47
1:A:533:SER:HB2	1:A:578:HIS:CD2	2.50	0.46
1:A:298:ILE:HD13	1:A:380:ILE:HD11	1.96	0.46
1:A:138:GLN:OE1	1:A:138:GLN:CA	2.64	0.46
1:A:229:ASN:ND2	1:A:232:GLU:CG	2.76	0.46
1:A:355:SER:HB3	1:A:385:THR:CG2	2.45	0.46
1:A:283:ARG:CD	1:A:283:ARG:N	2.58	0.46
1:A:624:THR:C	1:A:626:HIS:H	2.24	0.46
1:A:206:GLU:HA	1:A:301:VAL:HA	1.98	0.46
1:A:480:THR:HG22	1:A:482:GLY:H	1.78	0.46
1:A:612:ASN:HA	1:A:615:LEU:HG	1.97	0.46
1:A:106:PHE:CB	1:A:151:GLU:HG2	2.25	0.46
1:A:430:THR:HB	1:A:458:LEU:HD22	1.97	0.46
1:A:640:LYS:NZ	1:A:640:LYS:CB	2.78	0.46
1:A:160:THR:O	1:A:160:THR:CG2	2.63	0.46
1:A:310:PRO:O	1:A:312:VAL:HG13	2.16	0.46
1:A:218:ARG:NH2	1:A:235:ASP:OD1	2.49	0.46
1:A:313:ILE:HB	1:A:316:TYR:HB2	1.97	0.46
1:A:452:LEU:HD23	1:A:452:LEU:C	2.40	0.46
1:A:506:TRP:CD2	1:A:597:ARG:CD	2.99	0.46
1:A:181:ASN:HA	1:A:182:PRO:HD3	1.76	0.45
1:A:533:SER:C	1:A:535:GLU:H	2.23	0.45
1:A:110:LEU:HB3	1:A:112:THR:HG23	1.99	0.45
1:A:143:MSE:HE2	1:A:148:ILE:HG21	1.99	0.45
1:A:207:VAL:HG11	1:A:303:TYR:HB3	1.99	0.45
1:A:531:ASP:CB	1:A:541:TYR:CE1	2.95	0.45
1:A:154:PRO:HG3	1:A:620:LEU:HD13	1.98	0.45
1:A:482:GLY:O	1:A:483:GLU:HB2	2.16	0.45
1:A:502:VAL:CG2	1:A:516:GLY:HA3	2.47	0.45
1:A:248:LEU:HB3	1:A:376:PHE:CD1	2.52	0.45
1:A:243:GLN:O	1:A:247:ASP:OD2	2.33	0.45
1:A:309:GLN:CB	1:A:310:PRO:HD3	2.46	0.45
1:A:159:GLN:HG2	1:A:281:GLU:OE2	2.17	0.45
1:A:192:PHE:O	1:A:195:ILE:HG12	2.17	0.45
1:A:220:GLY:O	1:A:229:ASN:HB2	2.16	0.45
1:A:562:ILE:O	1:A:566:VAL:HG23	2.16	0.45
1:A:226:GLU:OE1	1:A:291:GLU:HG3	2.17	0.45
1:A:330:ARG:HG2	1:A:513:HIS:CE1	2.52	0.45
1:A:416:MSE:O	1:A:419:VAL:HG23	2.17	0.45
1:A:535:GLU:C	1:A:537:GLY:N	2.71	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:558:LEU:O	1:A:562:ILE:HG13	2.17	0.45
1:A:188:LEU:HD23	1:A:188:LEU:HA	1.87	0.44
1:A:209:TRP:O	1:A:213:GLU:CG	2.63	0.44
1:A:409:ASP:CG	1:A:409:ASP:O	2.60	0.44
1:A:191:ILE:O	1:A:194:THR:HG23	2.16	0.44
1:A:302:MSE:HE3	1:A:377:ILE:HG12	1.99	0.44
1:A:403:ARG:HG3	1:A:404:GLY:N	2.32	0.44
1:A:95:GLU:C	1:A:97:ASN:N	2.75	0.44
1:A:336:PHE:HB3	1:A:455:GLU:OE1	2.18	0.44
1:A:151:GLU:H	1:A:151:GLU:HG3	1.31	0.44
1:A:184:LEU:O	1:A:188:LEU:HG	2.18	0.44
1:A:289:TYR:HA	1:A:290:PRO:HD3	1.84	0.44
1:A:348:ALA:O	1:A:349:TRP:C	2.61	0.44
1:A:647:VAL:C	1:A:649:PRO:CD	2.90	0.44
1:A:282:LYS:HG2	1:A:283:ARG:HD3	2.00	0.44
1:A:153:LEU:O	1:A:155:ILE:N	2.50	0.44
1:A:229:ASN:ND2	1:A:230:ILE:O	2.50	0.44
1:A:541:TYR:O	1:A:545:VAL:HG23	2.17	0.44
1:A:100:ILE:HG23	1:A:101:LEU:N	2.32	0.44
1:A:126:ASP:N	1:A:130:ARG:O	2.28	0.43
1:A:263:LYS:HD3	1:A:265:GLU:OE2	2.18	0.43
1:A:184:LEU:HD22	1:A:188:LEU:HG	2.00	0.43
1:A:115:MSE:HE2	1:A:115:MSE:HB3	1.90	0.43
1:A:195:ILE:HD12	1:A:314:PRO:CG	2.48	0.43
1:A:649:PRO:HG3	1:A:661:TYR:CD2	2.54	0.43
1:A:290:PRO:HB2	1:A:295:ARG:HG3	2.01	0.43
1:A:309:GLN:CB	1:A:310:PRO:CD	2.96	0.43
1:A:342:VAL:HB	1:A:452:LEU:CD2	2.49	0.43
1:A:528:THR:O	1:A:530:LEU:CD1	2.67	0.43
1:A:355:SER:CB	1:A:385:THR:HG23	2.48	0.43
1:A:106:PHE:HE1	1:A:122:SER:HA	1.83	0.43
1:A:218:ARG:HH22	1:A:235:ASP:CG	2.27	0.43
1:A:105:ARG:HG3	1:A:122:SER:HB3	2.00	0.43
1:A:678:GLN:O	1:A:678:GLN:CG	2.44	0.43
1:A:403:ARG:NH2	1:A:413:GLY:HA3	2.34	0.43
1:A:533:SER:CB	1:A:578:HIS:CD2	3.01	0.43
1:A:646:LEU:O	1:A:649:PRO:HD2	2.18	0.43
1:A:286:VAL:HG23	1:A:286:VAL:O	2.18	0.43
1:A:387:VAL:HA	1:A:388:PRO:HD3	1.89	0.43
1:A:499:HIS:HE1	1:A:553:TYR:OH	2.02	0.43
1:A:531:ASP:CB	1:A:541:TYR:CZ	3.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:208:THR:C	1:A:210:GLU:H	2.27	0.42
1:A:277:ASP:OD1	1:A:277:ASP:C	2.62	0.42
1:A:648:LYS:HZ1	1:A:653:ILE:HG22	1.84	0.42
1:A:147:GLY:HA3	1:A:624:THR:HG23	2.01	0.42
1:A:289:TYR:HB3	1:A:406:GLY:HA3	2.01	0.42
1:A:508:ASP:C	1:A:508:ASP:OD2	2.61	0.42
1:A:142:ILE:O	1:A:145:SER:HB2	2.18	0.42
1:A:363:GLU:OE2	1:A:363:GLU:HA	2.18	0.42
1:A:277:ASP:OD1	1:A:279:VAL:CG1	2.59	0.42
1:A:320:THR:O	1:A:321:PRO:C	2.63	0.42
1:A:247:ASP:O	1:A:252:ARG:HB2	2.20	0.42
1:A:398:ILE:O	1:A:398:ILE:HG22	2.20	0.42
1:A:594:VAL:HG13	1:A:668:THR:HB	2.02	0.42
1:A:206:GLU:HB2	1:A:373:TRP:CZ2	2.55	0.42
1:A:283:ARG:HD3	1:A:283:ARG:N	2.24	0.42
1:A:650:ASP:OD2	1:A:653:ILE:HD13	2.19	0.42
1:A:648:LYS:HA	1:A:648:LYS:NZ	2.33	0.41
1:A:143:MSE:CE	1:A:148:ILE:HG21	2.50	0.41
1:A:154:PRO:HG2	1:A:620:LEU:CD1	2.50	0.41
1:A:175:LYS:CD	1:A:351:THR:HG23	2.32	0.41
1:A:274:GLN:HE21	1:A:620:LEU:HD11	1.85	0.41
1:A:355:SER:O	1:A:359:GLN:HG3	2.20	0.41
1:A:325:ILE:HD11	1:A:496:PHE:CE2	2.55	0.41
1:A:640:LYS:O	1:A:659:HIS:CD2	2.74	0.41
1:A:263:LYS:HG3	1:A:287:ILE:HD11	2.02	0.41
1:A:124:GLN:HG3	1:A:134:ILE:HG12	2.02	0.41
1:A:200:LEU:C	1:A:203:THR:HG23	2.45	0.41
1:A:100:ILE:CG2	1:A:101:LEU:N	2.84	0.41
1:A:403:ARG:NH2	1:A:413:GLY:C	2.79	0.41
1:A:623:TRP:CD1	1:A:623:TRP:C	2.97	0.41
1:A:154:PRO:HB3	1:A:270:SER:HA	2.02	0.41
1:A:165:PHE:CE1	1:A:262:PRO:HD3	2.56	0.41
1:A:204:TYR:HB2	1:A:369:TYR:HA	2.03	0.40
1:A:177:GLU:OE1	1:A:354:THR:HG23	2.21	0.40
1:A:504:VAL:HG11	1:A:602:LEU:CD2	2.52	0.40
1:A:202:HIS:HE1	1:A:374:HIS:HE1	1.70	0.40
1:A:348:ALA:O	1:A:352:GLN:NE2	2.55	0.40
1:A:352:GLN:HE21	1:A:352:GLN:HB2	1.64	0.40
1:A:220:GLY:O	1:A:221:ALA:O	2.39	0.40
1:A:531:ASP:CG	1:A:533:SER:H	2.29	0.40
1:A:536:ARG:HA	1:A:536:ARG:NH1	2.29	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:GLU:OE1	1:A:128:GLU:OE1[4_675]	2.14	0.06

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	586/609 (96%)	525 (90%)	47 (8%)	14 (2%)	4	24

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	221	ALA
1	A	309	GLN
1	A	219	LYS
1	A	277	ASP
1	A	401	GLY
1	A	537	GLY
1	A	154	PRO
1	A	237	GLU
1	A	535	GLU
1	A	646	LEU
1	A	625	LYS
1	A	96	HIS
1	A	293	LYS
1	A	648	LYS

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	516/524 (98%)	467 (90%)	49 (10%)	8	31

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	125	LEU
1	A	151	GLU
1	A	154	PRO
1	A	184	LEU
1	A	189	LEU
1	A	194	THR
1	A	199	THR
1	A	227	LYS
1	A	229	ASN
1	A	236	SER
1	A	244	LEU
1	A	248	LEU
1	A	254	ILE
1	A	274	GLN
1	A	279	VAL
1	A	283	ARG
1	A	290	PRO
1	A	293	LYS
1	A	301	VAL
1	A	317	GLU
1	A	322	LEU
1	A	340	VAL
1	A	345	ASP
1	A	351	THR
1	A	352	GLN
1	A	419	VAL
1	A	435	LYS
1	A	445	VAL
1	A	452	LEU
1	A	480	THR
1	A	531	ASP
1	A	536	ARG
1	A	538	THR
1	A	539	THR
1	A	550	LEU

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Mol	Chain	Res	Type
1	A	554	SER
1	A	558	LEU
1	A	599	LEU
1	A	600	SER
1	A	602	LEU
1	A	610	LEU
1	A	640	LYS
1	A	641	GLU
1	A	646	LEU
1	A	647	VAL
1	A	648	LYS
1	A	674	TYR
1	A	675	GLU
1	A	678	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	159	GLN
1	A	202	HIS
1	A	229	ASN
1	A	274	GLN
1	A	352	GLN
1	A	365	GLN
1	A	418	ASN
1	A	468	GLN
1	A	499	HIS
1	A	513	HIS
1	A	556	ASN
1	A	598	ASN
1	A	670	GLN
1	A	673	HIS

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand 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	GTP	A	2030	-	33,34,34	0.80	1 (3%)	50,54,54	0.96	1 (2%)

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	2030	-	-	6/22/38/38	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	2030	GTP	PA-O3A	2.88	1.62	1.59

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2030	GTP	C2'-C1'-N9	4.36	125.40	113.25

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	2030	GTP	PB-O3B-PG-O3G

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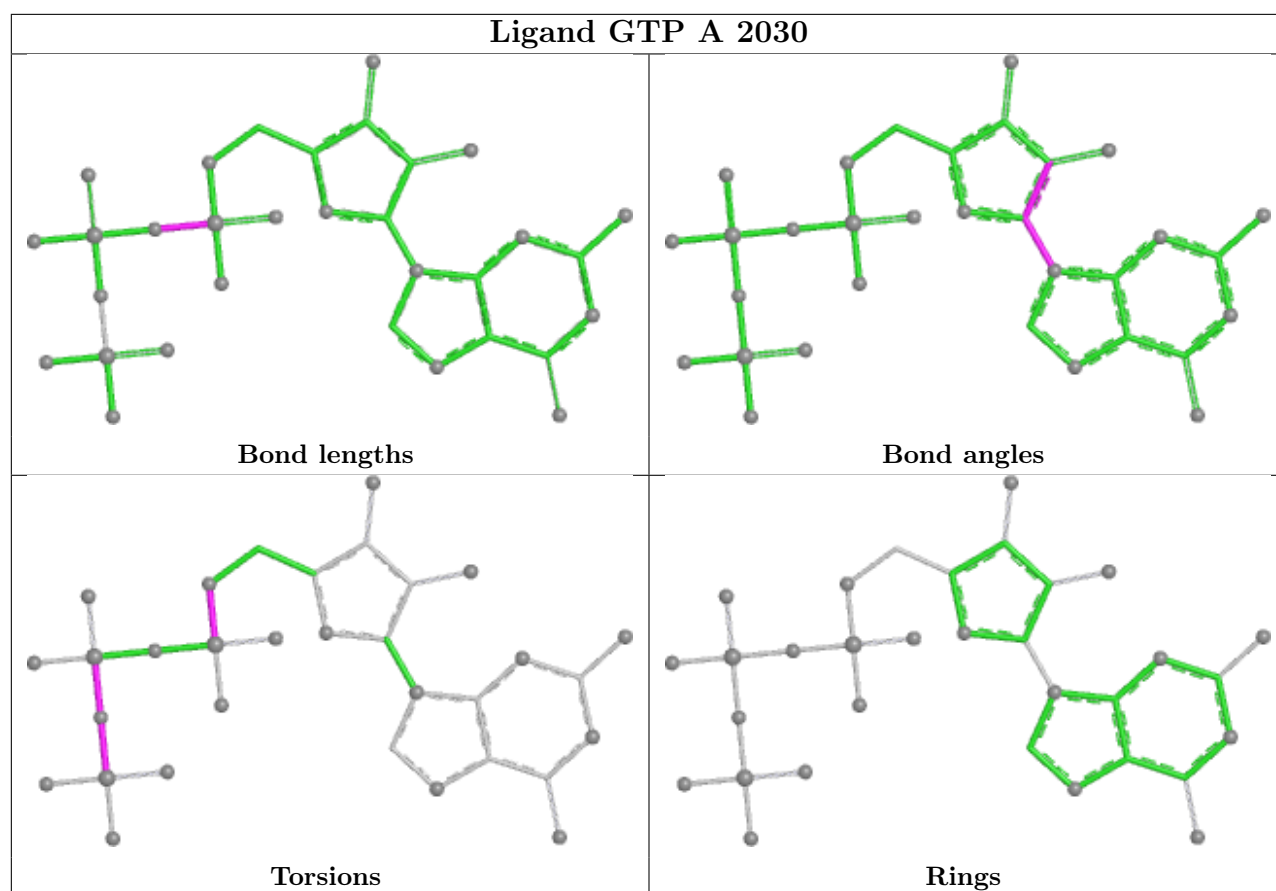
Mol	Chain	Res	Type	Atoms
2	A	2030	GTP	C5'-O5'-PA-O1A
2	A	2030	GTP	PG-O3B-PB-O2B
2	A	2030	GTP	PB-O3B-PG-O1G
2	A	2030	GTP	PB-O3B-PG-O2G
2	A	2030	GTP	PG-O3B-PB-O1B

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2030	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 ⓘ

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	576/609 (94%)	0.54	61 (10%) 11 6	6, 30, 69, 86	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	531	ASP	4.9
1	A	674	TYR	4.2
1	A	533	SER	3.9
1	A	671	GLY	3.9
1	A	676	GLN	3.7
1	A	130	ARG	3.5
1	A	532	SER	3.4
1	A	537	GLY	3.4
1	A	538	THR	3.3
1	A	227	LYS	3.3
1	A	216	VAL	3.2
1	A	253	LYS	3.1
1	A	672	LYS	3.1
1	A	536	ARG	3.1
1	A	539	THR	3.0
1	A	159	GLN	3.0
1	A	535	GLU	3.0
1	A	678	GLN	2.9
1	A	572	GLU	2.9
1	A	131	LYS	2.8
1	A	371	LYS	2.8
1	A	221	ALA	2.8
1	A	218	ARG	2.8
1	A	530	LEU	2.8
1	A	673	HIS	2.7
1	A	147	GLY	2.7
1	A	215	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	675	GLU	2.5
1	A	102	LYS	2.5
1	A	239	HIS	2.5
1	A	209	TRP	2.5
1	A	292	ALA	2.4
1	A	576	SER	2.4
1	A	401	GLY	2.4
1	A	529	ARG	2.4
1	A	435	LYS	2.4
1	A	534	GLY	2.3
1	A	264	ASN	2.3
1	A	667	PHE	2.3
1	A	234	LEU	2.3
1	A	482	GLY	2.3
1	A	210	GLU	2.3
1	A	211	GLN	2.3
1	A	129	GLY	2.2
1	A	291	GLU	2.2
1	A	251	GLY	2.2
1	A	626	HIS	2.2
1	A	92	VAL	2.2
1	A	123	GLU	2.2
1	A	208	THR	2.1
1	A	507	SER	2.1
1	A	240	LEU	2.1
1	A	310	PRO	2.1
1	A	126	ASP	2.1
1	A	283	ARG	2.1
1	A	309	GLN	2.1
1	A	331	LYS	2.0
1	A	133	ASN	2.0
1	A	625	LYS	2.0
1	A	157	ARG	2.0
1	A	109	ASN	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

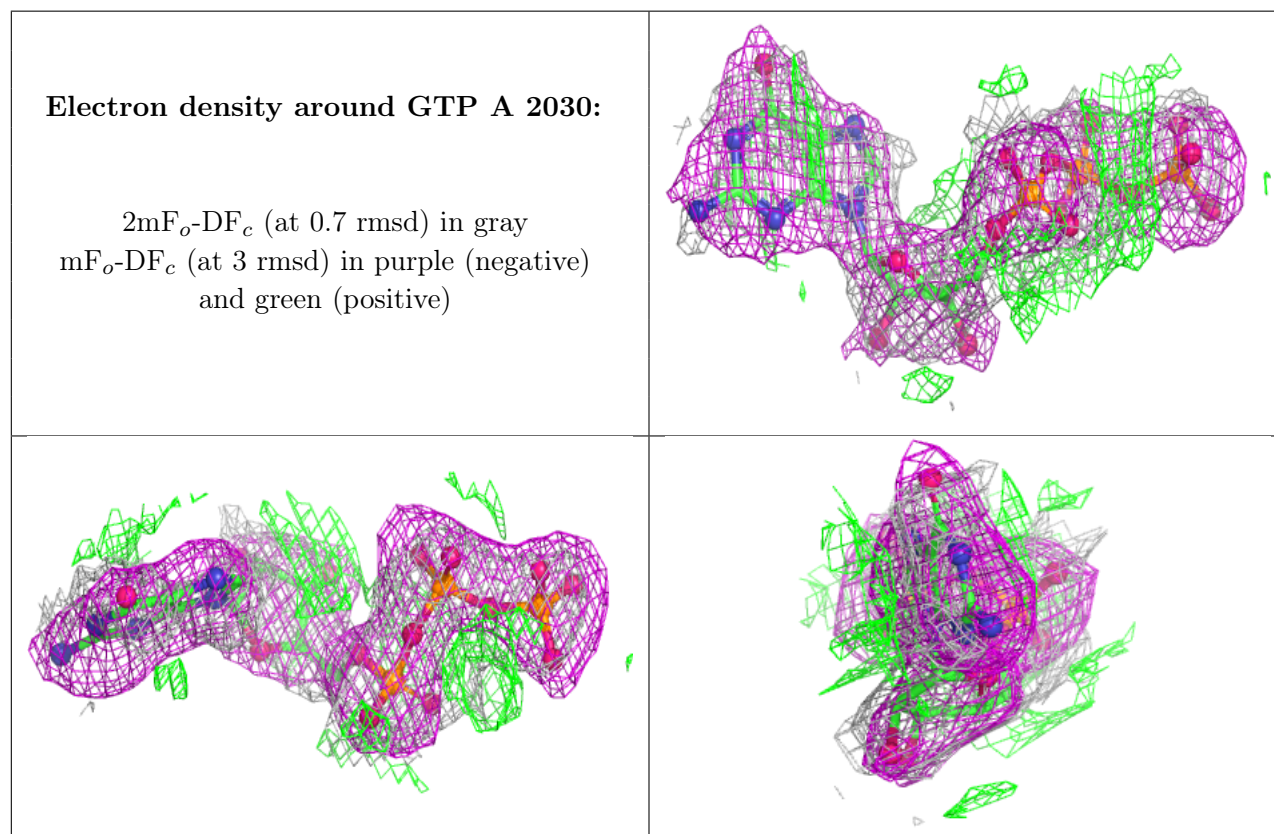
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
2	GTP	A	2030	32/32	0.66	0.23	6,6,6,6	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.