



Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2020 – 10:19 pm BST

PDB ID : 1Z5W
Title : Crystal Structure of gamma-tubulin bound to GTP
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Deposited on : 2005-03-20
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 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 : 2.11
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) : 2.11

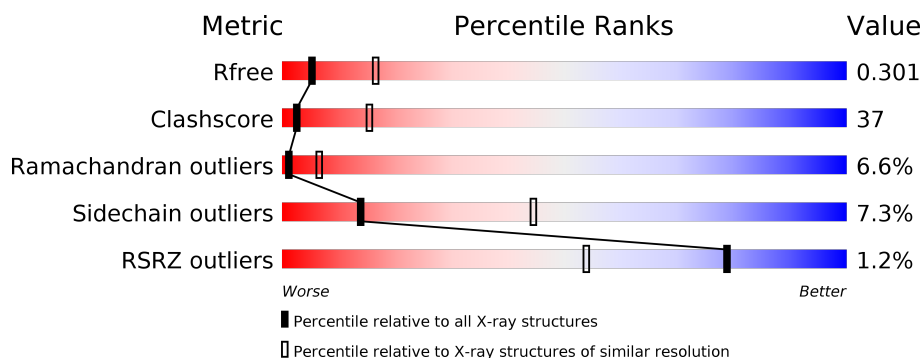
1 Overall quality at a glance

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}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (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 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	474	<div> <div></div> <div> <div></div> <div>39%</div> <div>41%</div> <div>7%</div> <div>14%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3110 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 Tubulin gamma-1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	409	3077	1951	524	588	14	7	0	0

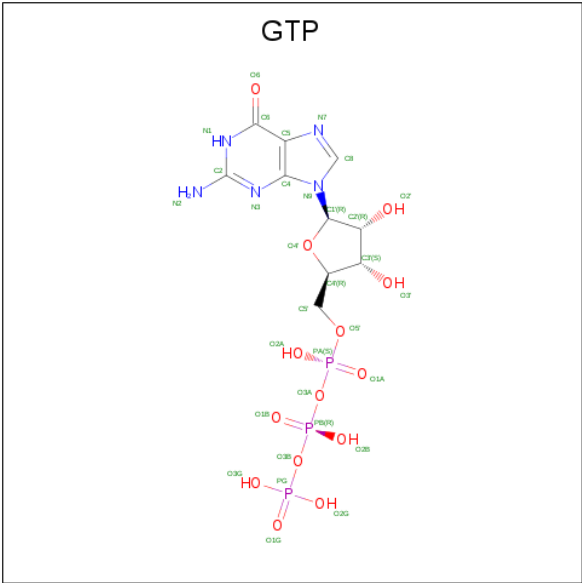
There are 25 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	450	VAL	-	CLONING ARTIFACT	UNP P23258
A	451	ASP	-	CLONING ARTIFACT	UNP P23258
A	452	VAL	-	CLONING ARTIFACT	UNP P23258
A	453	ASP	-	CLONING ARTIFACT	UNP P23258
A	454	GLY	-	CLONING ARTIFACT	UNP P23258
A	455	GLY	-	CLONING ARTIFACT	UNP P23258
A	456	GLU	-	CLONING ARTIFACT	UNP P23258
A	457	GLN	-	CLONING ARTIFACT	UNP P23258
A	458	LYS	-	CLONING ARTIFACT	UNP P23258
A	459	LEU	-	CLONING ARTIFACT	UNP P23258
A	460	ILE	-	CLONING ARTIFACT	UNP P23258
A	461	SER	-	CLONING ARTIFACT	UNP P23258
A	462	GLU	-	CLONING ARTIFACT	UNP P23258
A	463	GLU	-	CLONING ARTIFACT	UNP P23258
A	464	ASP	-	CLONING ARTIFACT	UNP P23258
A	465	LEU	-	CLONING ARTIFACT	UNP P23258
A	466	LEU	-	CLONING ARTIFACT	UNP P23258
A	467	LEU	-	CLONING ARTIFACT	UNP P23258
A	468	GLU	-	CLONING ARTIFACT	UNP P23258
A	469	HIS	-	EXPRESSION TAG	UNP P23258
A	470	HIS	-	EXPRESSION TAG	UNP P23258
A	471	HIS	-	EXPRESSION TAG	UNP P23258
A	472	HIS	-	EXPRESSION TAG	UNP P23258
A	473	HIS	-	EXPRESSION TAG	UNP P23258
A	474	HIS	-	EXPRESSION TAG	UNP P23258

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).

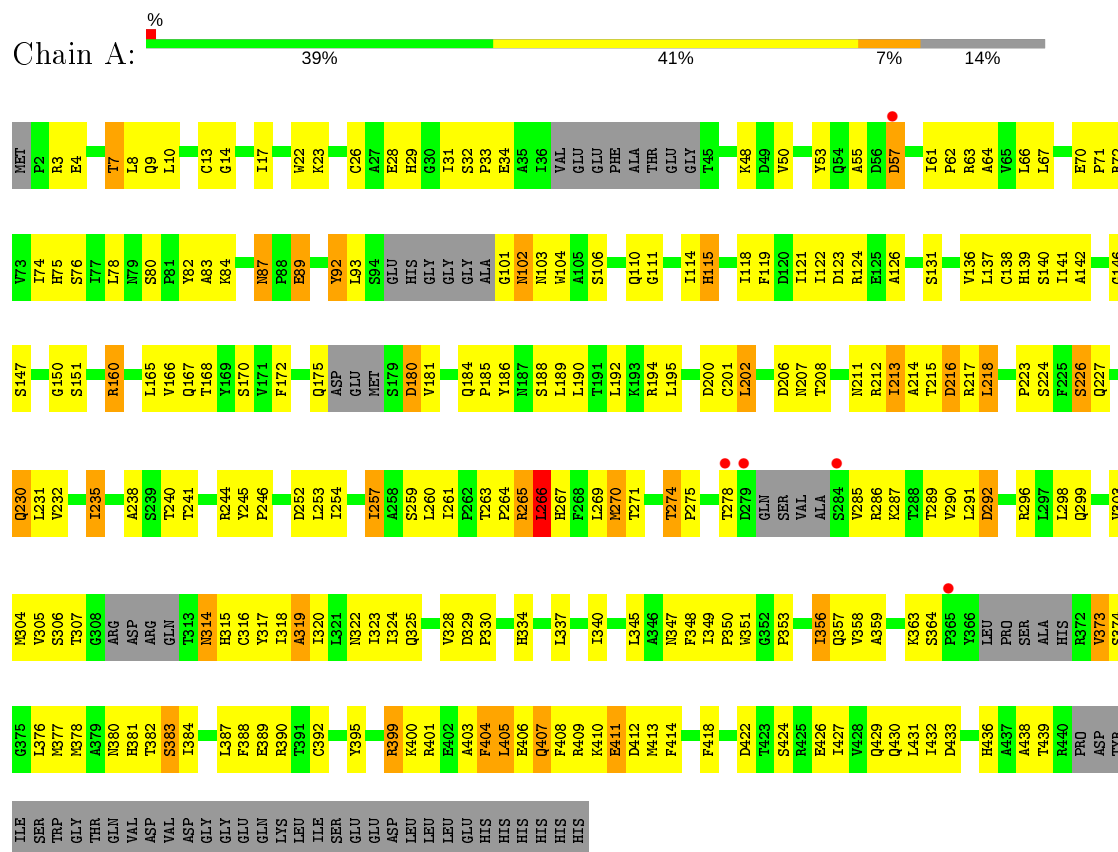


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

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 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: Tubulin gamma-1 chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.19Å 76.29Å 64.81Å 90.00° 102.31° 90.00°	Depositor
Resolution (Å)	29.91 – 3.00 29.91 – 2.50	Depositor EDS
% Data completeness (in resolution range)	85.8 (29.91-3.00) 80.9 (29.91-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.00 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.247 , 0.309 0.242 , 0.301	Depositor DCC
R_{free} test set	1580 reflections (10.05%)	wwPDB-VP
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.683	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 52.4	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	0.88	EDS
Total number of atoms	3110	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.55% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.42	0/3133	0.70	0/4266

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	92	TYR	Sidechain

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3077	0	2888	216	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	32	0	12	5	0
All	All	3110	0	2900	219	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 37.

All (219) 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:84:LYS:HE2	1:A:84:LYS:HA	1.43	0.98
1:A:266:LEU:HD13	1:A:432:ILE:HD13	1.46	0.97
1:A:165:LEU:HD13	1:A:254:ILE:HG22	1.46	0.96
1:A:373:VAL:HG12	1:A:374:SER:H	1.31	0.95
1:A:195:LEU:HD22	1:A:202:LEU:HD11	1.49	0.92
1:A:290:VAL:HG21	1:A:328:VAL:HG13	1.52	0.91
1:A:409:ARG:HA	1:A:413:MET:HB3	1.51	0.90
1:A:141:ILE:HD11	1:A:170:SER:HB3	1.54	0.89
1:A:70:GLU:HG2	1:A:71:PRO:HD2	1.54	0.86
1:A:296:ARG:HA	1:A:299:GLN:HG3	1.55	0.86
1:A:141:ILE:HG22	1:A:188:SER:HA	1.58	0.82
1:A:103:ASN:ND2	1:A:407:GLN:HG3	1.95	0.82
1:A:259:SER:CB	1:A:357:GLN:HE22	1.94	0.80
1:A:28:GLU:HB3	1:A:245:TYR:OH	1.80	0.79
1:A:238:ALA:O	1:A:241:THR:HG23	1.83	0.79
1:A:3:ARG:HD3	1:A:131:SER:OG	1.83	0.78
1:A:87:ASN:C	1:A:87:ASN:HD22	1.87	0.78
1:A:87:ASN:ND2	1:A:89:GLU:OE1	2.16	0.78
1:A:151:SER:HB3	1:A:194:ARG:HG3	1.67	0.77
1:A:238:ALA:HB3	1:A:376:LEU:HD13	1.68	0.76
1:A:274:THR:HB	1:A:275:PRO:CD	2.16	0.76
1:A:426:GLU:O	1:A:430:GLN:HG2	1.85	0.75
1:A:160:ARG:HG3	1:A:160:ARG:HH11	1.53	0.74
1:A:257:ILE:HG13	1:A:261:ILE:HD12	1.69	0.73
1:A:292:ASP:O	1:A:296:ARG:HD2	1.89	0.73
1:A:406:GLU:OE1	1:A:406:GLU:HA	1.89	0.72
1:A:389:GLU:HA	1:A:429:GLN:HE22	1.54	0.72
1:A:167:GLN:HG2	1:A:168:THR:N	2.03	0.72
1:A:382:THR:HG23	1:A:432:ILE:O	1.89	0.72
1:A:314:ASN:HB2	1:A:436:HIS:CD2	2.26	0.71
1:A:218:LEU:HD12	1:A:218:LEU:O	1.92	0.70
1:A:17:ILE:HD11	1:A:138:CYS:HB3	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:CYS:HB2	1:A:257:ILE:HD12	1.74	0.69
1:A:29:HIS:HB3	1:A:48:LYS:HE2	1.73	0.69
1:A:8:LEU:HB2	1:A:137:LEU:HD12	1.73	0.69
1:A:240:THR:O	1:A:244:ARG:HG3	1.92	0.69
1:A:32:SER:C	1:A:34:GLU:H	1.97	0.68
1:A:406:GLU:O	1:A:407:GLN:HB3	1.94	0.68
1:A:181:VAL:HB	1:A:184:GLN:HG3	1.75	0.68
1:A:264:PRO:O	1:A:266:LEU:N	2.26	0.67
1:A:274:THR:HB	1:A:275:PRO:HD3	1.77	0.67
1:A:373:VAL:HG12	1:A:374:SER:N	2.07	0.66
1:A:350:PRO:HG2	1:A:438:ALA:HA	1.77	0.66
1:A:84:LYS:HE2	1:A:84:LYS:CA	2.23	0.66
1:A:121:ILE:HG23	1:A:124:ARG:HH22	1.60	0.66
1:A:3:ARG:O	1:A:50:VAL:HG13	1.94	0.66
1:A:103:ASN:HD22	1:A:407:GLN:HG3	1.60	0.66
1:A:285:VAL:HG12	1:A:286:ARG:H	1.61	0.65
1:A:9:GLN:NE2	1:A:66:LEU:HD22	2.12	0.65
1:A:224:SER:OG	1:A:227:GLN:HG3	1.97	0.65
1:A:285:VAL:HG12	1:A:286:ARG:N	2.13	0.64
1:A:389:GLU:CA	1:A:429:GLN:HE22	2.11	0.64
1:A:10:LEU:HD11	1:A:137:LEU:HD11	1.80	0.64
1:A:32:SER:O	1:A:34:GLU:N	2.30	0.64
1:A:238:ALA:HB3	1:A:376:LEU:CD1	2.27	0.63
1:A:101:GLY:O	1:A:102:ASN:HB3	1.99	0.62
1:A:78:LEU:O	1:A:83:ALA:HB2	1.98	0.62
1:A:160:ARG:HG3	1:A:160:ARG:NH1	2.10	0.62
1:A:184:GLN:N	1:A:185:PRO:HD2	2.13	0.62
1:A:10:LEU:O	1:A:14:GLY:HA3	1.99	0.62
1:A:166:VAL:N	1:A:200:ASP:OD2	2.30	0.62
3:A:2466:GTP:H5''	3:A:2466:GTP:H8	1.64	0.62
1:A:254:ILE:HA	1:A:257:ILE:CG2	2.30	0.61
1:A:253:LEU:O	1:A:257:ILE:HG22	2.01	0.61
1:A:270:MET:CE	1:A:303:VAL:HG11	2.30	0.61
1:A:101:GLY:HA3	1:A:106:SER:OG	2.02	0.60
1:A:181:VAL:HG21	1:A:184:GLN:HE21	1.65	0.60
1:A:325:GLN:OE1	1:A:363:LYS:HG2	2.01	0.60
1:A:211:ASN:O	1:A:212:ARG:C	2.40	0.59
1:A:215:THR:HG22	1:A:216:ASP:OD1	2.03	0.59
1:A:151:SER:CB	1:A:194:ARG:HG3	2.32	0.58
1:A:31:ILE:HG22	1:A:32:SER:O	2.03	0.58
1:A:340:ILE:HG23	1:A:345:LEU:HB2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:PHE:HA	1:A:122:ILE:HB	1.85	0.58
1:A:252:ASP:HB3	1:A:254:ILE:HG12	1.85	0.58
1:A:121:ILE:HG23	1:A:124:ARG:NH2	2.18	0.58
1:A:382:THR:O	1:A:384:ILE:N	2.37	0.58
1:A:404:PHE:C	1:A:406:GLU:H	2.06	0.57
1:A:412:ASP:HA	1:A:414:PHE:CE1	2.39	0.57
1:A:74:ILE:HD12	1:A:93:LEU:HD21	1.85	0.57
1:A:424:SER:O	1:A:427:ILE:HB	2.05	0.57
1:A:316:CYS:HB3	1:A:347:ASN:ND2	2.20	0.56
1:A:403:ALA:O	1:A:405:LEU:N	2.38	0.56
1:A:266:LEU:HD21	1:A:431:LEU:HB3	1.87	0.56
1:A:382:THR:C	1:A:384:ILE:H	2.08	0.56
3:A:2466:GTP:H5"	3:A:2466:GTP:C8	2.40	0.56
1:A:324:ILE:HD13	1:A:358:VAL:HG13	1.87	0.56
1:A:181:VAL:HB	1:A:184:GLN:CG	2.35	0.56
1:A:350:PRO:HD2	1:A:351:TRP:CD1	2.41	0.56
1:A:232:VAL:O	1:A:235:ILE:HG23	2.06	0.55
1:A:70:GLU:HG2	1:A:71:PRO:CD	2.32	0.55
1:A:289:THR:OG1	1:A:291:LEU:HG	2.05	0.55
1:A:186:TYR:OH	1:A:405:LEU:HD21	2.07	0.54
1:A:261:ILE:HD13	1:A:267:HIS:HB2	1.89	0.54
1:A:111:GLY:O	1:A:115:HIS:N	2.41	0.54
1:A:141:ILE:O	1:A:147:SER:HB3	2.08	0.54
1:A:408:PHE:C	1:A:410:LYS:H	2.11	0.54
1:A:307:THR:HG21	1:A:383:SER:HB2	1.90	0.53
1:A:325:GLN:HB3	1:A:363:LYS:CG	2.39	0.53
1:A:376:LEU:C	1:A:376:LEU:HD23	2.29	0.53
1:A:257:ILE:CG1	1:A:261:ILE:HD12	2.38	0.53
1:A:266:LEU:HD23	1:A:266:LEU:N	2.23	0.53
1:A:101:GLY:O	1:A:102:ASN:CB	2.58	0.52
1:A:92:TYR:HE2	1:A:114:ILE:HD13	1.74	0.52
1:A:404:PHE:C	1:A:406:GLU:N	2.63	0.52
1:A:392:CYS:O	1:A:395:TYR:N	2.43	0.52
1:A:175:GLN:N	1:A:208:THR:OG1	2.40	0.52
1:A:114:ILE:O	1:A:115:HIS:C	2.48	0.52
1:A:87:ASN:C	1:A:87:ASN:ND2	2.61	0.52
1:A:314:ASN:HB2	1:A:436:HIS:HD2	1.72	0.51
1:A:214:ALA:O	1:A:218:LEU:HB3	2.10	0.51
1:A:305:VAL:HG13	1:A:387:LEU:HD11	1.91	0.51
1:A:53:TYR:HD2	1:A:61:ILE:HD11	1.75	0.51
1:A:325:GLN:HB3	1:A:363:LYS:HG2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:LEU:HD11	1:A:422:ASP:HA	1.91	0.51
1:A:238:ALA:CB	1:A:376:LEU:HD13	2.40	0.51
1:A:389:GLU:N	1:A:429:GLN:HE22	2.09	0.51
1:A:433:ASP:O	1:A:436:HIS:ND1	2.32	0.51
1:A:377:MET:HG2	1:A:378:MET:N	2.26	0.51
1:A:317:TYR:CG	1:A:381:HIS:HD2	2.29	0.51
1:A:186:TYR:HE2	1:A:395:TYR:HE1	1.60	0.50
1:A:167:GLN:CG	1:A:168:THR:N	2.74	0.50
1:A:323:ILE:HG12	1:A:359:ALA:HB3	1.93	0.50
1:A:76:SER:O	1:A:80:SER:HB3	2.12	0.49
1:A:13:CYS:SG	1:A:140:SER:HB3	2.51	0.49
1:A:119:PHE:CD2	1:A:160:ARG:HD2	2.48	0.49
1:A:67:LEU:CD2	1:A:92:TYR:HB3	2.42	0.49
1:A:114:ILE:HG22	1:A:118:ILE:HG13	1.95	0.49
1:A:404:PHE:O	1:A:406:GLU:N	2.45	0.49
1:A:320:ILE:HD12	1:A:348:PHE:CE2	2.48	0.49
1:A:426:GLU:OE2	1:A:430:GLN:NE2	2.44	0.49
1:A:74:ILE:HD12	1:A:93:LEU:CD2	2.42	0.49
1:A:315:HIS:ND1	1:A:382:THR:CG2	2.77	0.48
1:A:334:HIS:HA	1:A:337:LEU:HD12	1.95	0.48
1:A:388:PHE:HB2	1:A:429:GLN:CD	2.33	0.48
1:A:350:PRO:HD2	1:A:351:TRP:HD1	1.79	0.48
1:A:413:MET:SD	1:A:418:PHE:CE2	3.07	0.48
1:A:4:GLU:HG2	1:A:63:ARG:CZ	2.44	0.48
1:A:66:LEU:N	1:A:66:LEU:HD23	2.29	0.48
1:A:87:ASN:HD21	1:A:89:GLU:CD	2.16	0.47
1:A:390:ARG:HG3	1:A:390:ARG:HH11	1.79	0.47
1:A:104:TRP:HD1	1:A:147:SER:HB2	1.79	0.47
1:A:240:THR:OG1	1:A:244:ARG:NH1	2.46	0.47
1:A:320:ILE:HD12	1:A:348:PHE:CZ	2.49	0.47
1:A:78:LEU:O	1:A:83:ALA:CB	2.62	0.47
1:A:213:ILE:O	1:A:214:ALA:C	2.53	0.47
1:A:304:MET:O	1:A:305:VAL:HG23	2.14	0.47
1:A:55:ALA:C	1:A:57:ASP:H	2.17	0.47
1:A:7:THR:HG23	1:A:136:VAL:HB	1.97	0.47
1:A:257:ILE:HG13	1:A:261:ILE:CD1	2.44	0.47
1:A:110:GLN:O	1:A:114:ILE:N	2.45	0.46
1:A:216:ASP:O	1:A:218:LEU:N	2.44	0.46
1:A:324:ILE:HD12	1:A:324:ILE:N	2.31	0.46
1:A:382:THR:C	1:A:384:ILE:N	2.69	0.46
1:A:395:TYR:O	1:A:399:ARG:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:GLN:N	1:A:185:PRO:CD	2.78	0.46
1:A:71:PRO:HB2	1:A:75:HIS:HD2	1.80	0.46
1:A:413:MET:SD	1:A:418:PHE:HE2	2.38	0.46
1:A:318:ILE:O	1:A:319:ALA:HB2	2.16	0.45
1:A:70:GLU:CG	1:A:71:PRO:HD2	2.35	0.45
1:A:215:THR:HG22	1:A:216:ASP:N	2.30	0.45
1:A:317:TYR:O	1:A:349:ILE:HG23	2.17	0.45
1:A:22:TRP:CZ2	1:A:64:ALA:HB2	2.52	0.45
1:A:195:LEU:HD22	1:A:202:LEU:CD1	2.35	0.45
1:A:373:VAL:CG1	1:A:374:SER:H	2.13	0.45
1:A:395:TYR:OH	1:A:418:PHE:HB3	2.17	0.45
1:A:13:CYS:HB2	3:A:2466:GTP:C8	2.51	0.45
1:A:254:ILE:HA	1:A:257:ILE:HG23	1.99	0.45
1:A:186:TYR:O	1:A:190:LEU:HG	2.17	0.45
1:A:23:LYS:HG3	1:A:82:TYR:CZ	2.52	0.45
1:A:358:VAL:HG12	1:A:359:ALA:N	2.31	0.44
1:A:244:ARG:C	1:A:246:PRO:HD3	2.38	0.44
1:A:410:LYS:O	1:A:411:GLU:HB3	2.17	0.44
1:A:61:ILE:HG13	1:A:61:ILE:O	2.17	0.44
1:A:146:GLY:O	1:A:150:GLY:HA3	2.17	0.44
1:A:260:LEU:HD13	1:A:269:LEU:HD13	2.00	0.44
1:A:399:ARG:C	1:A:401:ARG:H	2.21	0.44
1:A:223:PRO:HA	1:A:227:GLN:OE1	2.17	0.43
1:A:316:CYS:HB3	1:A:347:ASN:HD21	1.82	0.43
1:A:114:ILE:O	1:A:114:ILE:HG22	2.18	0.43
1:A:285:VAL:CG1	1:A:286:ARG:N	2.81	0.43
1:A:206:ASP:OD1	1:A:208:THR:HB	2.18	0.43
1:A:271:THR:HG22	1:A:378:MET:HG2	2.00	0.43
1:A:160:ARG:CG	1:A:160:ARG:HH11	2.23	0.43
1:A:167:GLN:HG3	1:A:201:CYS:HB3	2.01	0.43
1:A:175:GLN:H	1:A:208:THR:HG1	1.65	0.43
1:A:207:ASN:HD21	3:A:2466:GTP:H1'	1.84	0.43
1:A:270:MET:HE3	1:A:303:VAL:HG11	1.99	0.43
1:A:238:ALA:HA	1:A:241:THR:HG23	2.01	0.43
3:A:2466:GTP:O1B	3:A:2466:GTP:O2A	2.34	0.43
1:A:382:THR:O	1:A:432:ILE:HG21	2.19	0.43
1:A:10:LEU:CD2	1:A:67:LEU:HB2	2.49	0.43
1:A:406:GLU:O	1:A:407:GLN:CB	2.65	0.42
1:A:226:SER:OG	1:A:227:GLN:N	2.52	0.42
1:A:382:THR:O	1:A:432:ILE:CG2	2.67	0.42
1:A:186:TYR:CE2	1:A:395:TYR:HE1	2.38	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:TRP:CE3	1:A:62:PRO:HB3	2.54	0.42
1:A:103:ASN:HD21	1:A:407:GLN:HG3	1.80	0.42
1:A:10:LEU:HA	1:A:67:LEU:O	2.20	0.42
1:A:28:GLU:O	1:A:364:SER:HB2	2.19	0.42
1:A:266:LEU:HD21	1:A:431:LEU:CB	2.47	0.42
1:A:230:GLN:HG2	1:A:231:LEU:N	2.34	0.42
1:A:270:MET:HE3	1:A:303:VAL:CG1	2.50	0.42
1:A:403:ALA:C	1:A:405:LEU:H	2.23	0.42
1:A:123:ASP:O	1:A:126:ALA:HB3	2.20	0.41
1:A:168:THR:OG1	1:A:202:LEU:HD13	2.20	0.41
1:A:3:ARG:NH1	1:A:3:ARG:HG3	2.35	0.41
1:A:67:LEU:HD23	1:A:92:TYR:HB3	2.03	0.41
1:A:3:ARG:HD3	1:A:131:SER:O	2.19	0.41
1:A:322:ASN:ND2	1:A:356:ILE:HD11	2.36	0.41
1:A:265:ARG:O	1:A:267:HIS:N	2.53	0.41
1:A:172:PHE:O	1:A:207:ASN:HB2	2.21	0.41
1:A:63:ARG:HH11	1:A:63:ARG:HG2	1.86	0.41
1:A:315:HIS:CE1	1:A:382:THR:HG22	2.56	0.41
1:A:71:PRO:O	1:A:72:ARG:C	2.59	0.41
1:A:238:ALA:HA	1:A:241:THR:CG2	2.51	0.40
1:A:136:VAL:HA	1:A:167:GLN:O	2.21	0.40
1:A:192:LEU:HD23	1:A:192:LEU:HA	1.86	0.40
1:A:270:MET:CE	1:A:303:VAL:CG1	2.99	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	395/474 (83%)	314 (80%)	55 (14%)	26 (7%)	1 6

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	180	ASP
1	A	314	ASN
1	A	404	PHE
1	A	411	GLU
1	A	102	ASN
1	A	115	HIS
1	A	217	ARG
1	A	266	LEU
1	A	319	ALA
1	A	330	PRO
1	A	383	SER
1	A	407	GLN
1	A	33	PRO
1	A	142	ALA
1	A	216	ASP
1	A	265	ARG
1	A	278	THR
1	A	400	LYS
1	A	405	LEU
1	A	218	LEU
1	A	274	THR
1	A	287	LYS
1	A	353	PRO
1	A	373	VAL
1	A	213	ILE
1	A	329	ASP

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	317/421 (75%)	294 (93%)	23 (7%)	14 44

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	THR

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Mol	Chain	Res	Type
1	A	26	CYS
1	A	57	ASP
1	A	87	ASN
1	A	89	GLU
1	A	139	HIS
1	A	160	ARG
1	A	180	ASP
1	A	202	LEU
1	A	226	SER
1	A	230	GLN
1	A	235	ILE
1	A	257	ILE
1	A	263	THR
1	A	266	LEU
1	A	270	MET
1	A	292	ASP
1	A	298	LEU
1	A	306	SER
1	A	356	ILE
1	A	380	ASN
1	A	399	ARG
1	A	439	THR

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

Mol	Chain	Res	Type
1	A	16	GLN
1	A	75	HIS
1	A	87	ASN
1	A	102	ASN
1	A	110	GLN
1	A	158	ASN
1	A	184	GLN
1	A	198	ASN
1	A	267	HIS
1	A	347	ASN
1	A	357	GLN
1	A	380	ASN
1	A	381	HIS
1	A	429	GLN

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	GTP	A	2466	2	26,34,34	1.89	6 (23%)	33,54,54	2.77	13 (39%)

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	GTP	A	2466	2	-	4/18/38/38	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2466	GTP	C6-N1	5.41	1.42	1.33
3	A	2466	GTP	O4'-C1'	4.11	1.46	1.41
3	A	2466	GTP	C8-N7	-3.40	1.28	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2466	GTP	C4-N3	2.84	1.40	1.35
3	A	2466	GTP	C2-N1	2.26	1.39	1.35
3	A	2466	GTP	C2'-C1'	-2.26	1.50	1.53

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2466	GTP	C5-C6-N1	-7.90	112.62	123.43
3	A	2466	GTP	C6-N1-C2	5.89	125.28	115.93
3	A	2466	GTP	O5'-PA-O1A	-5.13	89.03	109.07
3	A	2466	GTP	O2A-PA-O5'	-5.07	84.21	107.75
3	A	2466	GTP	O2B-PB-O1B	4.53	134.63	112.24
3	A	2466	GTP	N2-C2-N3	3.26	123.10	117.79
3	A	2466	GTP	N3-C2-N1	-3.11	123.08	127.22
3	A	2466	GTP	O2A-PA-O1A	3.08	127.47	112.24
3	A	2466	GTP	O5'-C5'-C4'	2.88	118.91	108.99
3	A	2466	GTP	C3'-C2'-C1'	2.70	105.04	100.98
3	A	2466	GTP	C2-N3-C4	-2.63	112.35	115.36
3	A	2466	GTP	N2-C2-N1	-2.21	113.82	117.25
3	A	2466	GTP	O4'-C4'-C5'	-2.01	102.77	109.37

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	2466	GTP	C5'-O5'-PA-O1A
3	A	2466	GTP	PB-O3A-PA-O2A
3	A	2466	GTP	C5'-O5'-PA-O3A
3	A	2466	GTP	PA-O3A-PB-O2B

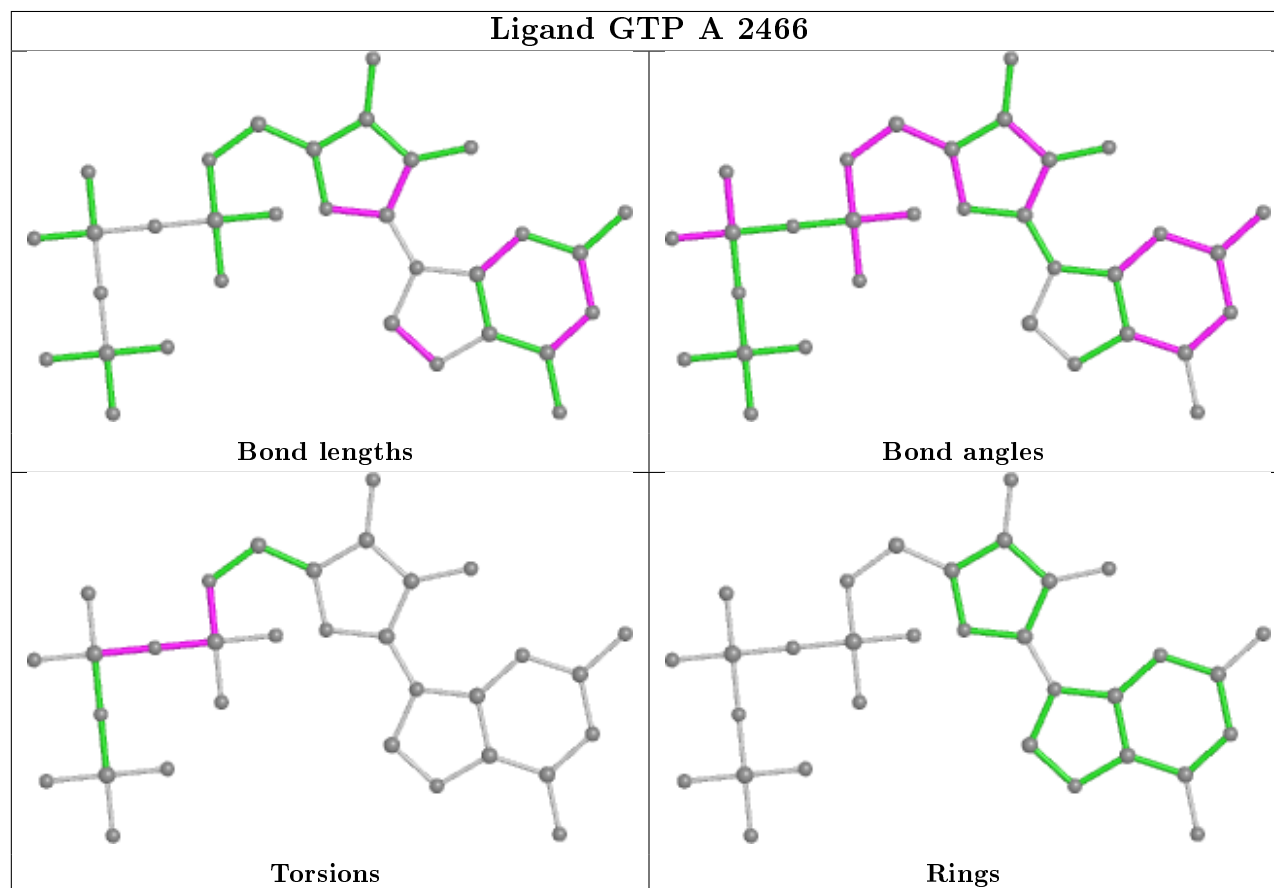
There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2466	GTP	5	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier.

Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	409/474 (86%)	-0.40	5 (1%) 79 54	21, 47, 83, 106	2 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	278	THR	2.7
1	A	284	SER	2.6
1	A	57	ASP	2.4
1	A	279	ASP	2.3
1	A	365	PRO	2.3

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](#)

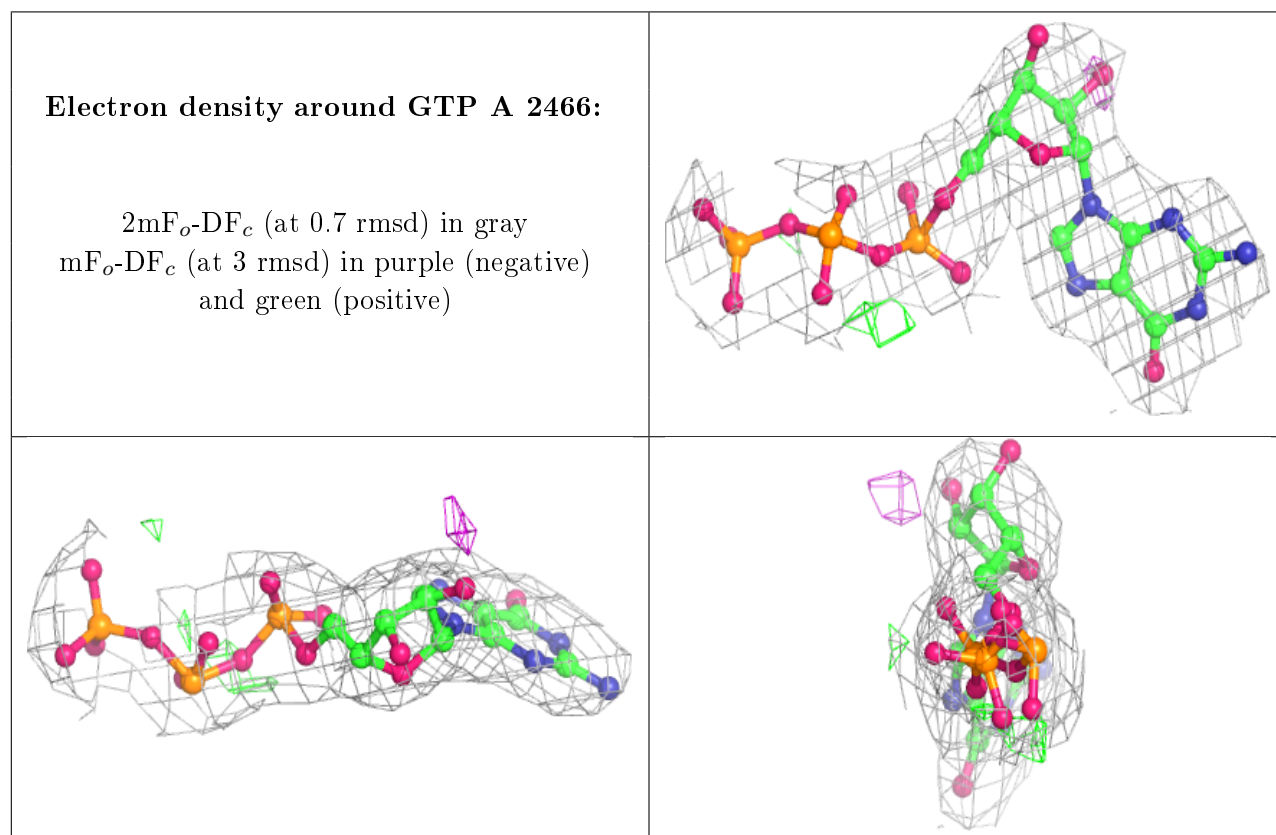
There are no carbohydrates in this entry.

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. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MG	A	475	1/1	0.94	0.24	53,53,53,53	0
3	GTP	A	2466	32/32	0.95	0.15	31,47,57,58	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.